

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017590.D
 Acq On : 06 Oct 2023 06:43
 Operator : MA/JU
 Sample : 04588-09
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHT7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017591.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.552	244	249	255	rBV2	1066528	1552787	8.79%	0.341%
2	5.263	365	370	373	rBV	1613243	2379356	13.47%	0.522%
3	5.399	389	393	403	rVB	1622519	2429415	13.75%	0.533%
4	5.705	440	445	449	rBV	978797	1578759	8.94%	0.346%
5	5.799	456	461	468	rBV	2241215	3557809	20.14%	0.781%
6	5.863	468	472	478	rVB	1170925	1790715	10.14%	0.393%
7	6.093	505	511	514	rBV	2410904	3992873	22.60%	0.876%
8	6.122	514	516	520	rVV	1951341	2924126	16.55%	0.642%
9	6.222	528	533	540	rVB2	2084228	3810345	21.57%	0.836%
10	6.310	540	548	551	rBV	1541272	2596684	14.70%	0.570%
11	6.393	551	562	566	rVV2	6277399	13276923	75.15%	2.914%
12	6.440	566	570	573	rVV2	1648628	2546283	14.41%	0.559%
13	6.493	573	579	584	rVV2	8205447	14517044	82.17%	3.186%
14	6.599	593	597	603	rVB	1503896	2097139	11.87%	0.460%
15	6.669	603	609	613	rBV3	1583502	2837253	16.06%	0.623%
16	6.710	613	616	618	rVV	1591117	2231512	12.63%	0.490%
17	6.734	618	620	625	rVB	2735399	3588897	20.31%	0.788%
18	6.793	625	630	632	rBV	1520839	2115480	11.97%	0.464%
19	6.834	632	637	641	rVV	5305447	8388615	47.48%	1.841%
20	6.887	644	646	651	rVB2	1943880	2462560	13.94%	0.540%
21	6.940	651	655	660	rVB4	1248630	2243061	12.70%	0.492%
22	7.005	660	666	672	rBV2	7507928	14256049	80.70%	3.129%
23	7.146	685	690	692	rVV	1036455	1626943	9.21%	0.357%
24	7.175	692	695	703	rVV2	2259964	4093369	23.17%	0.898%
25	7.246	703	707	711	rVV2	1842052	3303549	18.70%	0.725%
26	7.305	711	717	720	rVV3	2355454	5634450	31.89%	1.237%
27	7.346	720	724	728	rVV	5071690	8249757	46.70%	1.810%
28	7.387	728	731	734	rVV2	2064001	3305670	18.71%	0.725%
29	7.422	734	737	742	rVV	2540178	4040862	22.87%	0.887%
30	7.469	742	745	754	rVB6	739653	1664267	9.42%	0.365%
31	7.546	754	758	761	rBV	2088884	2841385	16.08%	0.624%
32	7.587	761	765	770	rVB4	2085899	3251945	18.41%	0.714%
33	7.640	770	774	777	rBV2	1577668	2307224	13.06%	0.506%
34	7.675	777	780	786	rVB3	1681709	2632229	14.90%	0.578%
35	7.769	786	796	798	rBV4	2903768	7016988	39.72%	1.540%
36	7.834	802	807	816	rVB	10676916	17666494	100.00%	3.877%

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017590.D
 Acq On : 06 Oct 2023 06:43
 Operator : MA/JU
 Sample : 04588-09
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHT7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Title : SVOA CALIBRATION

37	7.922	816	822	830	rBV5	2180653	5448056	30.84%	1.196%
38	8.046	837	843	850	rBV4	3926806	8080698	45.74%	1.773%
39	8.110	850	854	858	rVB2	2567352	3242887	18.36%	0.712%
40	8.169	858	864	868	rBV2	5668633	9552067	54.07%	2.096%

41	8.216	868	872	878	rVB3	2253460	3650614	20.66%	0.801%
42	8.310	879	888	893	rVB5	1916745	5125280	29.01%	1.125%
43	8.387	893	901	906	rBV3	5147198	10773926	60.99%	2.364%
44	8.446	906	911	919	rVB3	4301977	9018734	51.05%	1.979%
45	8.522	919	924	926	rBV2	2616878	3894812	22.05%	0.855%

46	8.552	926	929	934	rVB3	2305200	3139645	17.77%	0.689%
47	8.628	938	942	952	rVB2	2670902	5903981	33.42%	1.296%
48	8.734	954	960	965	rBV2	5261951	9891933	55.99%	2.171%
49	8.863	977	982	986	rBV	1172062	1896417	10.73%	0.416%
50	8.957	992	998	1000	rBV3	2126536	4123381	23.34%	0.905%

51	8.993	1000	1004	1006	rVV2	3743192	6528969	36.96%	1.433%
52	9.022	1006	1009	1015	rVV2	5004937	8346760	47.25%	1.832%
53	9.081	1015	1019	1022	rVV2	1904410	3339152	18.90%	0.733%
54	9.110	1022	1024	1028	rVB	1721213	2037559	11.53%	0.447%
55	9.204	1035	1040	1043	rBV3	1316347	2240534	12.68%	0.492%

56	9.246	1044	1047	1056	rVB7	2001662	3746630	21.21%	0.822%
57	9.346	1056	1064	1071	rVB3	2147720	5471931	30.97%	1.201%
58	9.510	1085	1092	1095	rBV3	2920592	5231368	29.61%	1.148%
59	9.551	1095	1099	1103	rVV	2381766	3762491	21.30%	0.826%
60	9.628	1103	1112	1119	rVB7	3810307	9188162	52.01%	2.016%

61	9.746	1124	1132	1136	rBV3	859678	2055877	11.64%	0.451%
62	9.816	1136	1144	1152	rBV4	3697551	8115997	45.94%	1.781%
63	9.893	1152	1157	1159	rVV	2476807	4037694	22.86%	0.886%
64	9.922	1159	1162	1170	rVB4	3654987	7101434	40.20%	1.558%
65	10.016	1170	1178	1186	rVB5	1684286	4404489	24.93%	0.967%

66	10.134	1187	1198	1203	rBV2	5014914	10005297	56.63%	2.196%
67	10.193	1203	1208	1214	rVB3	1382174	3076457	17.41%	0.675%
68	10.593	1272	1276	1280	rVV3	839958	1687448	9.55%	0.370%
69	10.657	1284	1287	1290	rVV2	957243	1650454	9.34%	0.362%
70	10.722	1294	1298	1302	rVV3	1956257	3364486	19.04%	0.738%

71	10.775	1302	1307	1313	rVV4	1525522	3934484	22.27%	0.863%
72	10.840	1313	1318	1325	rVB2	6222069	10711884	60.63%	2.351%
73	10.946	1325	1336	1343	rBV5	1170504	3665818	20.75%	0.804%
74	11.175	1371	1375	1381	rVB3	906452	1606783	9.10%	0.353%
75	11.410	1410	1415	1420	rBV4	2078609	3564211	20.17%	0.782%

76	11.528	1428	1435	1442	rBV8	991156	2316087	13.11%	0.508%
----	--------	------	------	------	------	--------	---------	--------	--------

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017590.D
 Acq On : 06 Oct 2023 06:43
 Operator : MA/JU
 Sample : 04588-09
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHT7

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Title : SVOA CALIBRATION

77	11.710	1461	1466	1471	rVV	6112092	8925910	50.52%	1.959%
78	11.975	1503	1511	1518	rBV5	1311317	2920585	16.53%	0.641%
79	12.334	1568	1572	1576	rVB	1684425	2238954	12.67%	0.491%
80	12.440	1585	1590	1600	rVB	1586779	2985735	16.90%	0.655%
81	12.657	1623	1627	1634	rBV2	1424331	2485311	14.07%	0.545%
82	12.757	1640	1644	1651	rBV5	1022500	2010904	11.38%	0.441%
83	12.828	1651	1656	1660	rVV2	1075503	2205301	12.48%	0.484%
84	13.075	1693	1698	1703	rBV	7362540	10136068	57.37%	2.224%
85	13.381	1748	1750	1758	rVV2	1211700	1688667	9.56%	0.371%
86	13.775	1811	1817	1818	rBV3	1294119	2007627	11.36%	0.441%
87	13.939	1840	1845	1850	rVV2	1426143	2417524	13.68%	0.531%
88	13.992	1850	1854	1856	rVV	1671037	2408647	13.63%	0.529%
89	14.028	1856	1860	1866	rVB2	6340977	9741233	55.14%	2.138%
90	14.316	1902	1909	1917	rBV6	971646	2776798	15.72%	0.609%
91	15.010	2021	2027	2032	rVB2	988777	1876722	10.62%	0.412%
92	15.092	2039	2041	2048	rVB2	1127690	1590692	9.00%	0.349%
93	15.292	2071	2075	2083	rVB3	974797	1825684	10.33%	0.401%
94	15.816	2158	2164	2170	rVB	4373730	6929773	39.23%	1.521%
95	16.304	2241	2247	2252	rBV	5277109	7839364	44.37%	1.720%
96	16.686	2305	2312	2319	rVB6	647215	1685332	9.54%	0.370%
97	17.139	2384	2389	2394	rBV	2373619	3494792	19.78%	0.767%
98	17.416	2431	2436	2440	rVV	1180465	1615204	9.14%	0.354%
99	21.551	3134	3139	3144	rBV	1557226	2011446	11.39%	0.441%
100	24.127	3570	3577	3588	rVB	1010597	2113581	11.96%	0.464%

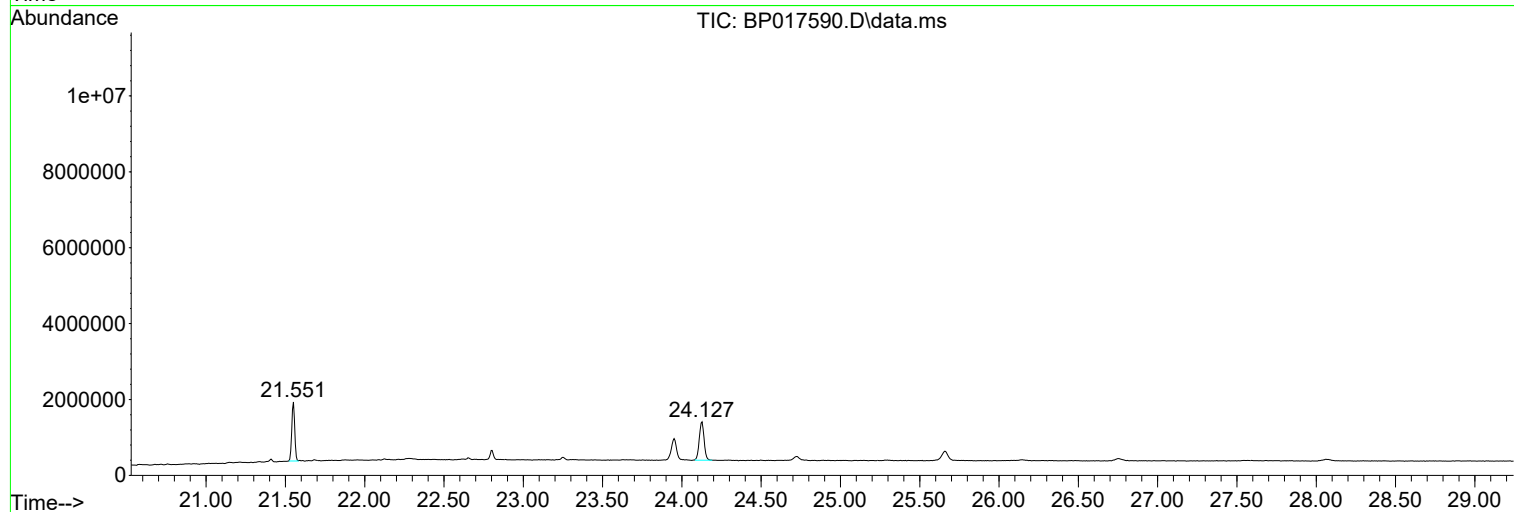
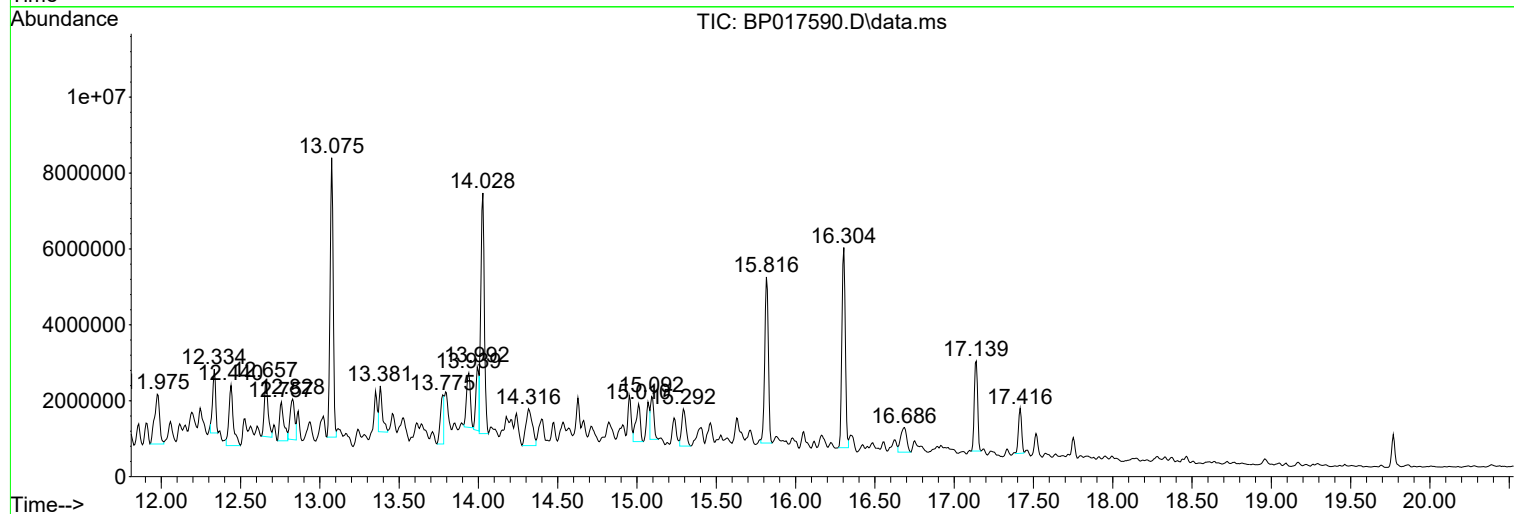
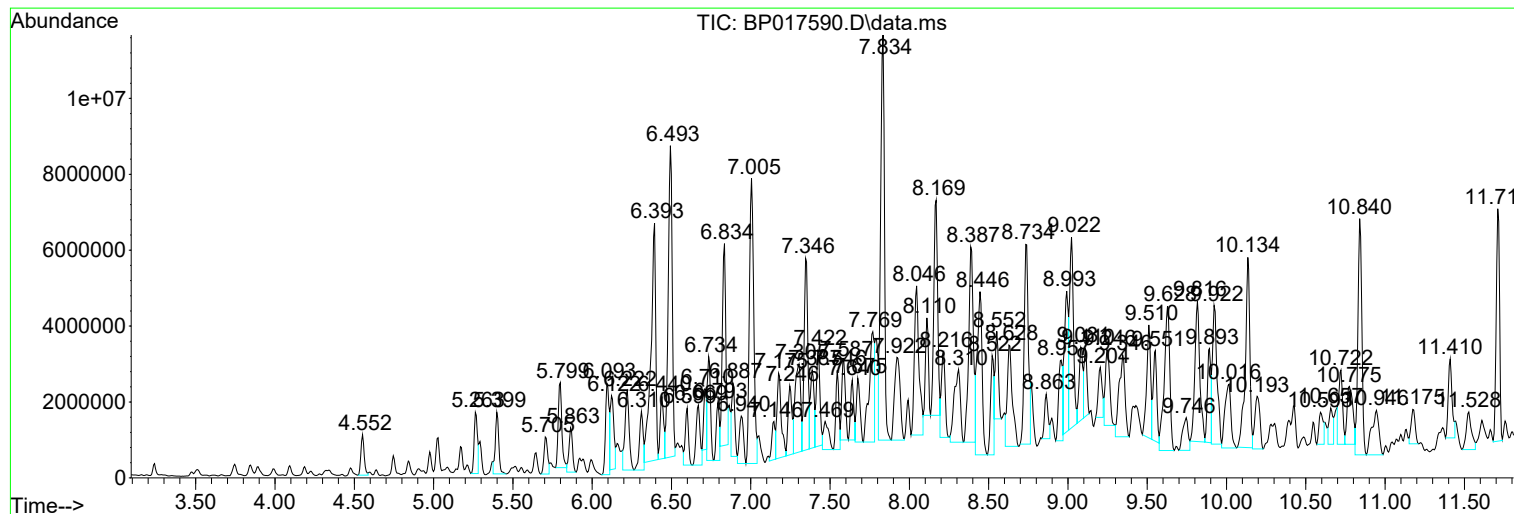
Sum of corrected areas: 455675558

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

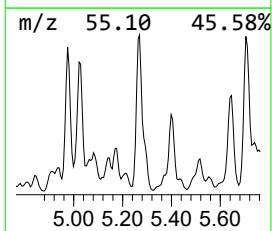
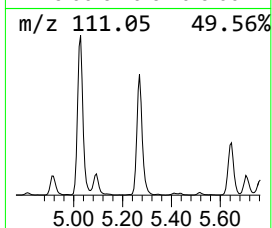
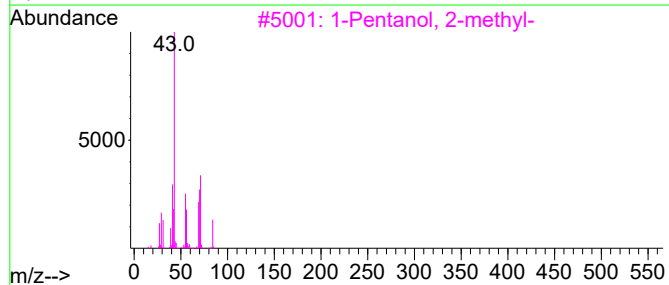
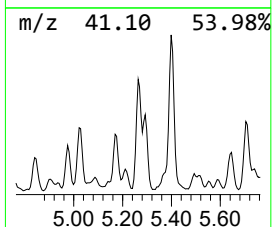
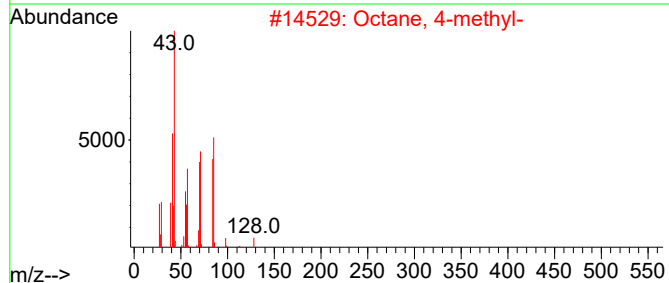
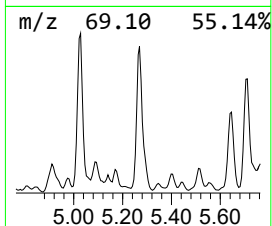
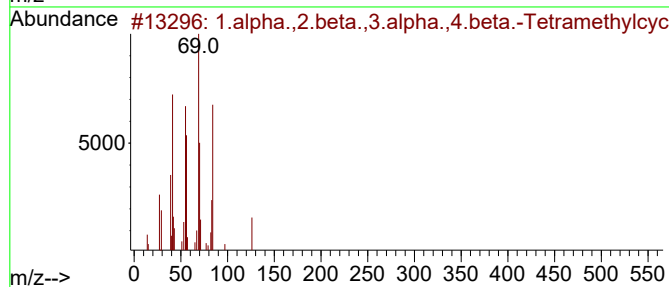
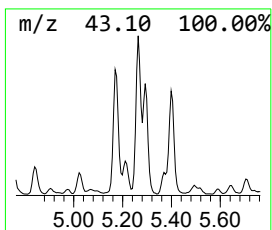
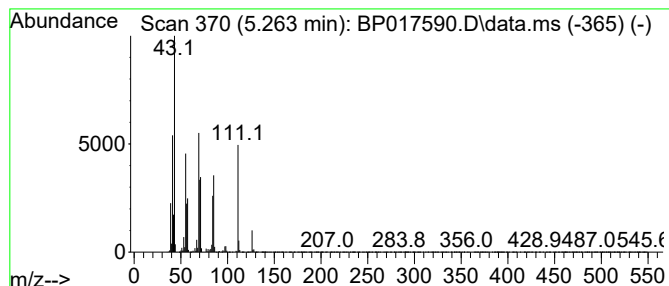
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 66

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.263	8.73 ng/ul	2379360	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1.alpha.,2.beta.,3.alpha.,4.beta...			126	C9H18	002532-67-4	43
2	Octane, 4-methyl-			128	C9H20	002216-34-4	43
3	1-Pentanol, 2-methyl-			102	C6H14O	000105-30-6	35
4	Heptane, 2,4-dimethyl-			128	C9H20	002213-23-2	27
5	Hexane, 2,3,5-trimethyl-			128	C9H20	001069-53-0	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

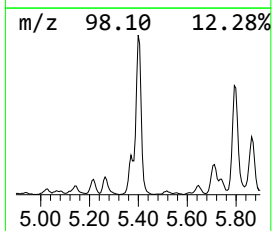
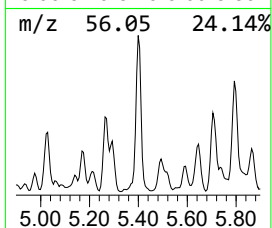
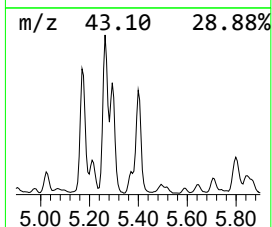
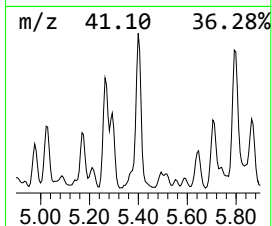
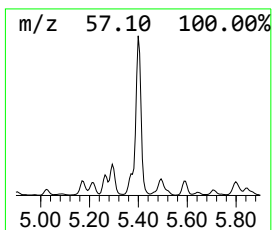
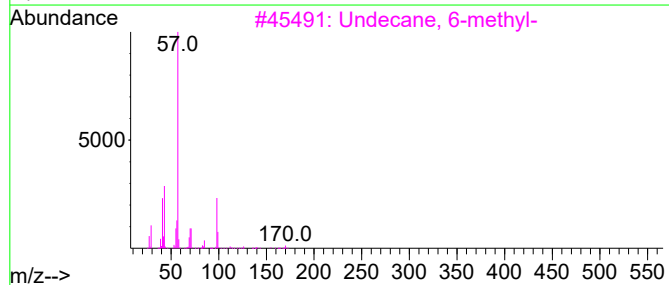
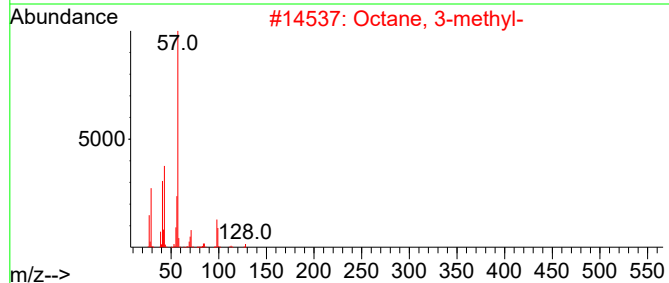
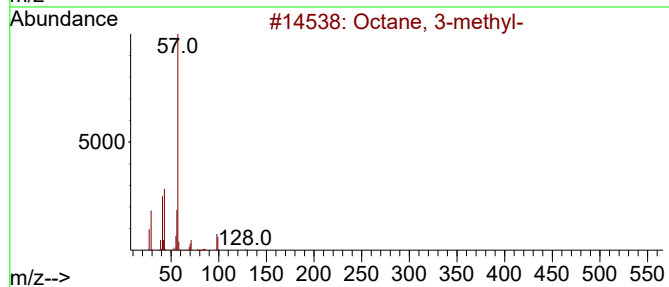
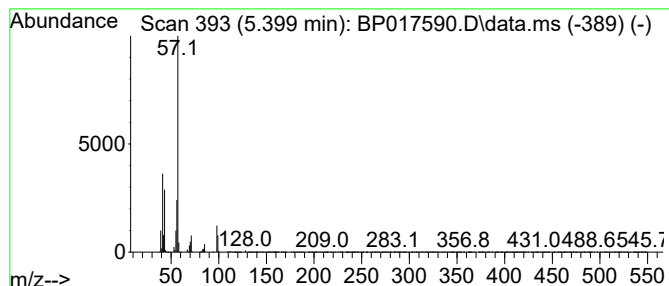
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 65

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.399	8.92 ng/ul	2429420	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-methyl-			128	C9H20	002216-33-3	90
2	Octane, 3-methyl-			128	C9H20	002216-33-3	83
3	Undecane, 6-methyl-			170	C12H26	017302-33-9	59
4	Undecane, 6-methyl-			170	C12H26	017302-33-9	59
5	Octane, 3-methyl-			128	C9H20	002216-33-3	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

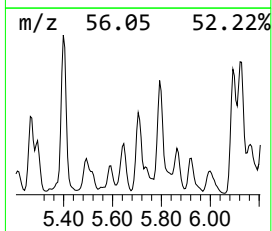
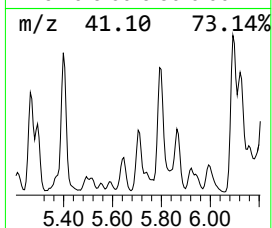
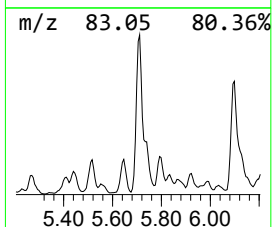
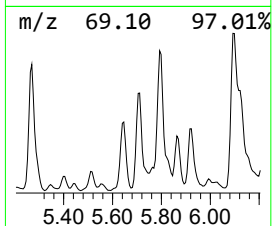
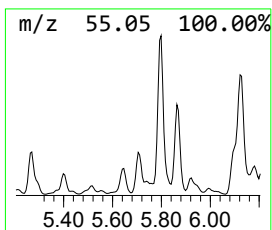
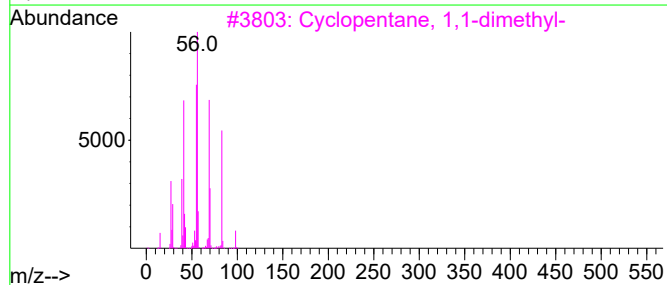
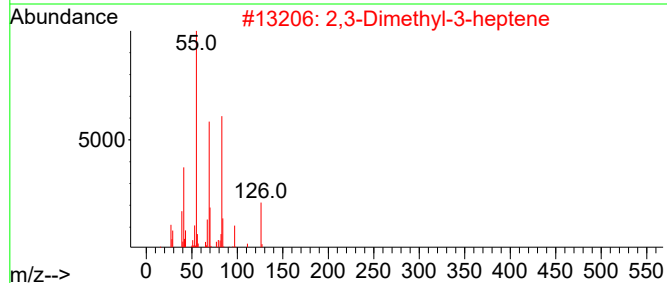
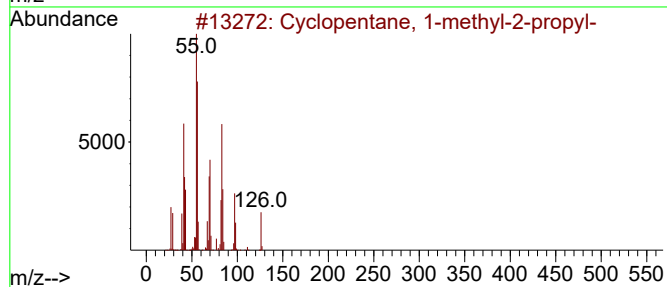
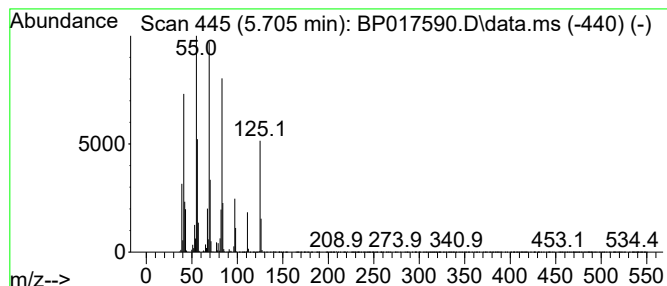
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 (DEL) Alkane: Cyclic5.705 Concentration Rank 79

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.705	5.80 ng/ul	1578760	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, 1-methyl-2-propyl-	126	C9H18	003728-57-2	60
2			2,3-Dimethyl-3-heptene	126	C9H18	1000113-49-3	53
3			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	50
4			Cyclohexane, 1,2,3-trimethyl-, (...)	126	C9H18	007667-55-2	49
5			Oxazole, 4-ethyl-2,5-dimethyl-	125	C7H11NO	030408-61-8	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

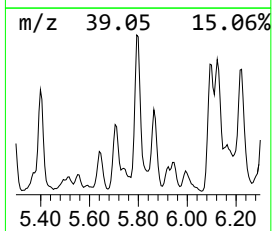
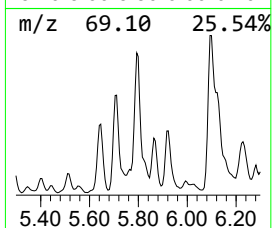
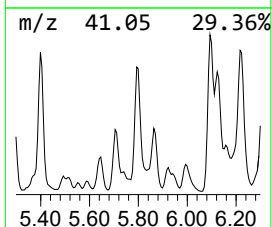
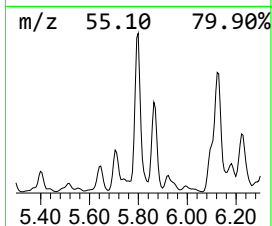
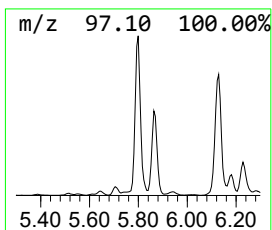
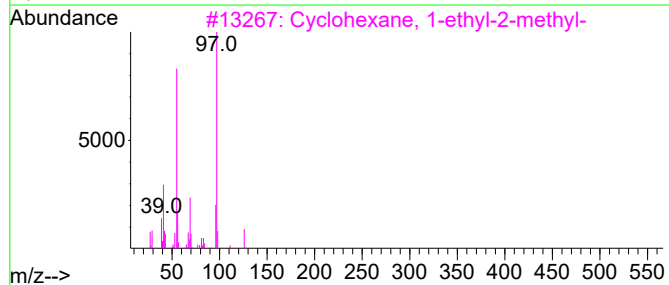
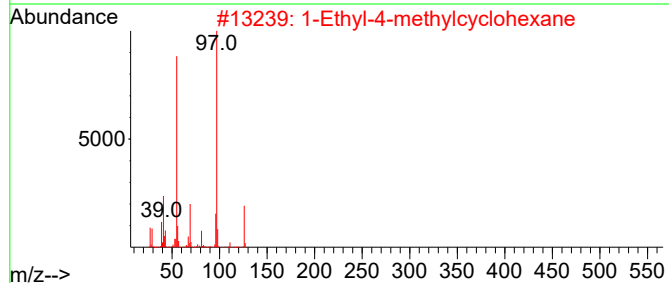
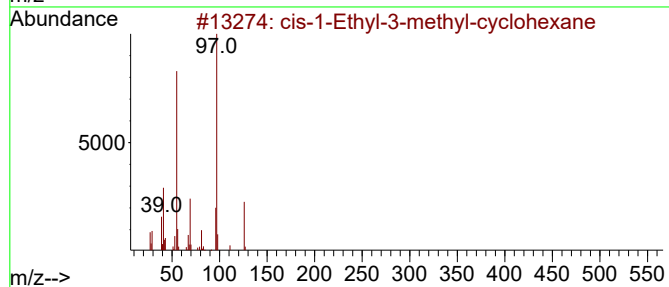
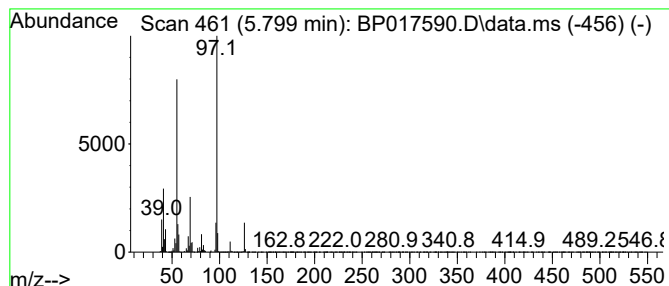
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 (DEL) Alkane: Cyclic5.799 Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.799	13.06 ng/ul	3557810	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	91
2			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	91
3			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	90
4			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	87
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

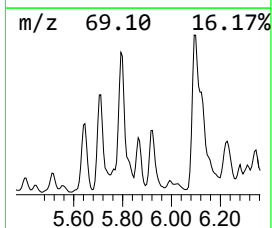
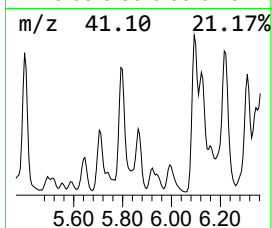
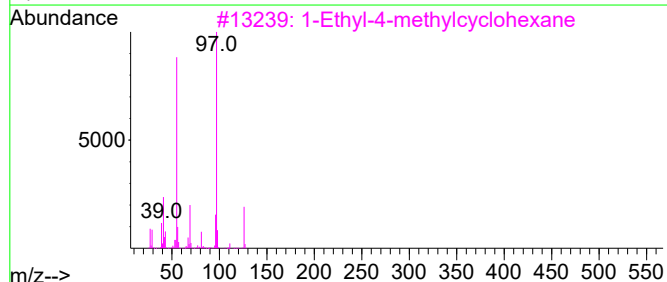
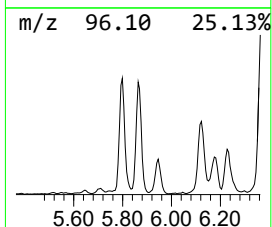
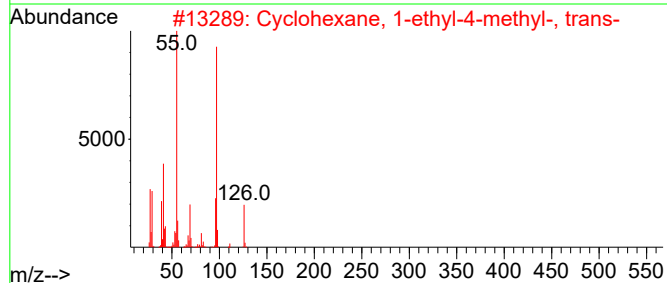
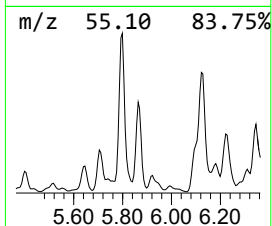
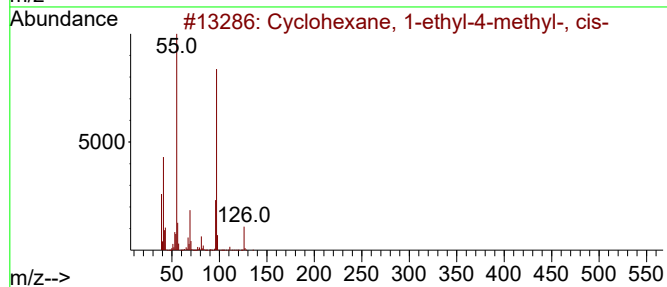
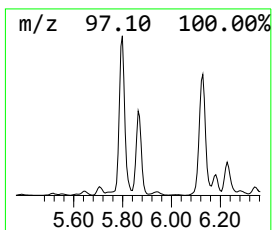
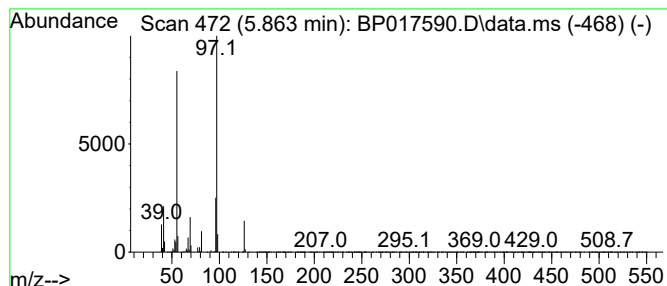
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 (DEL) Alkane: Cyclic5.863 Concentration Rank 77

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.863	6.57 ng/ul	1790720	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	91
2			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	006236-88-0	91
3			1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	91
4			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	91
5			Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

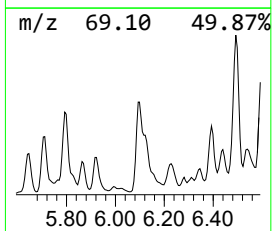
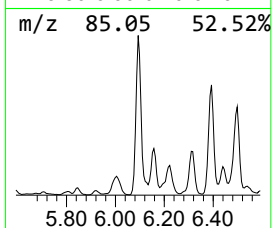
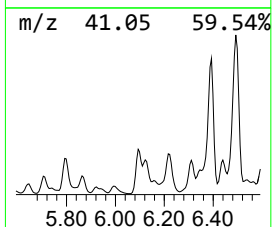
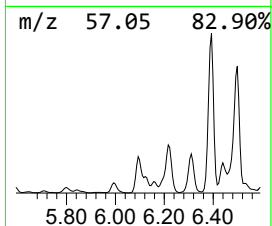
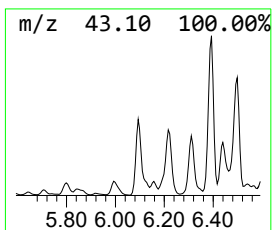
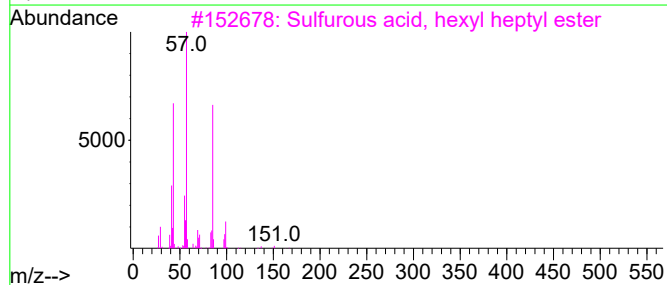
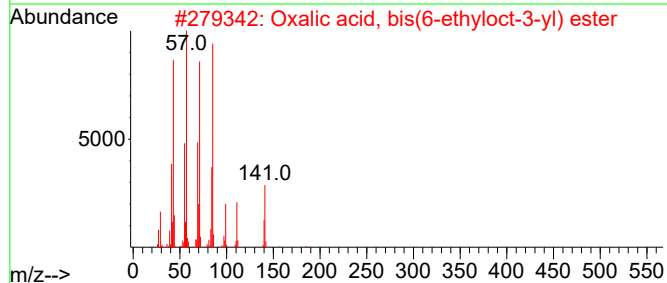
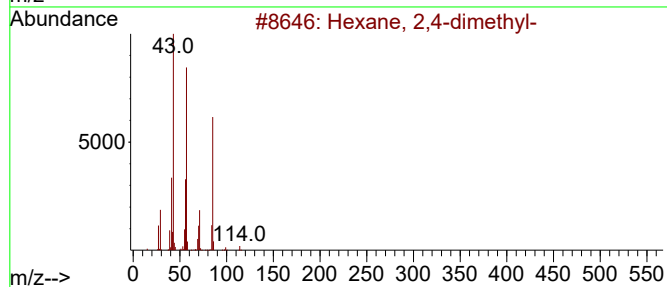
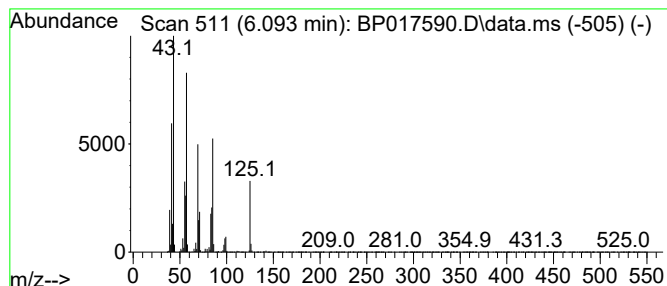
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.093	14.66 ng/ul	3992870	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	50
2	Oxalic acid, bis(6-ethyloct-3-yl...)	370	C22H42O4	1000309-34-6	50
3	Sulfurous acid, hexyl heptyl ester	264	C13H28O3S	1000309-12-9	47
4	Hexane, 2,4-dimethyl-	114	C8H18	000589-43-5	47
5	Ethanol, 2-(hexyloxy)-	146	C8H18O2	000112-25-4	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

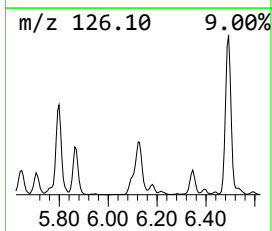
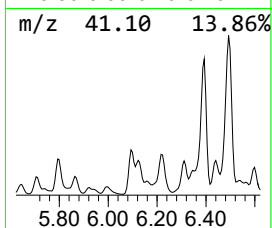
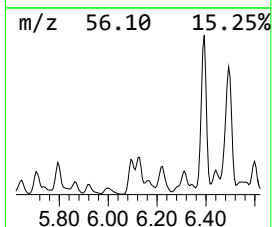
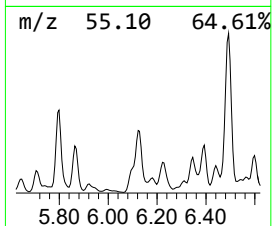
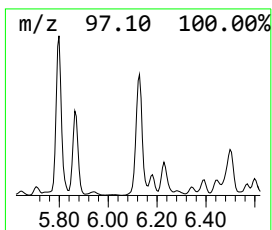
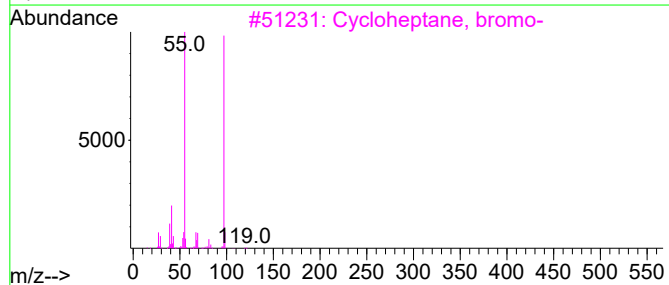
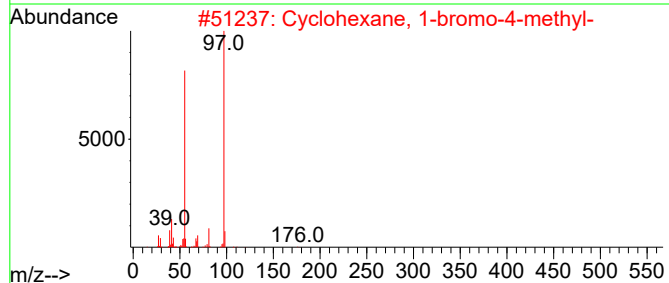
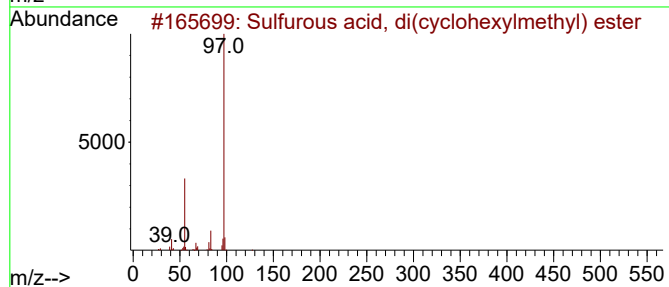
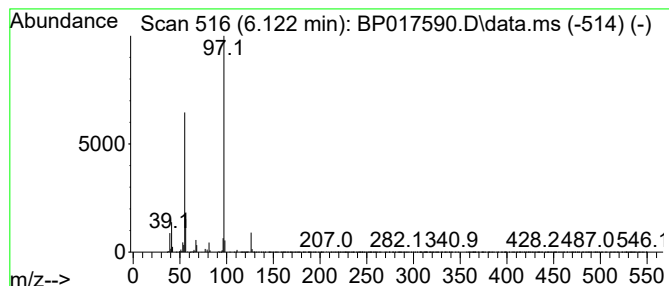
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 Sulfurous acid, di(cyclohex... Concentration Rank 57

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.122	10.73 ng/ul	2924130	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, di(cyclohexylmet...	274	C14H26O3S	1010309-22-7	72
2			Cyclohexane, 1-bromo-4-methyl-	176	C7H13Br	006294-40-2	64
3			Cycloheptane, bromo-	176	C7H13Br	002404-35-5	64
4			Cyclohexane, 1,3-dimethyl-	112	C8H16	000591-21-9	56
5			Succinic acid, cyclohexylmethyl ...	358	C17H20Cl2O4	1000390-15-3	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

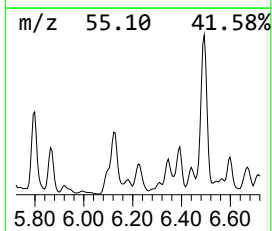
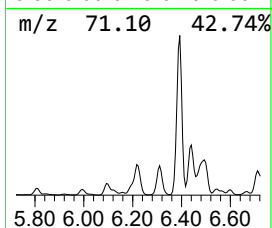
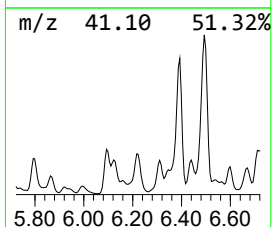
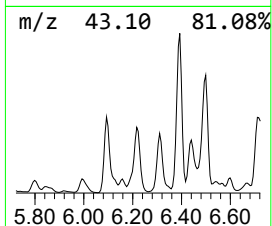
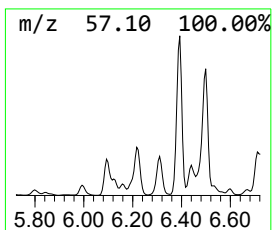
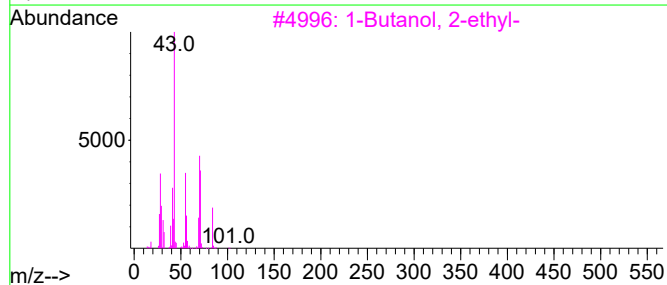
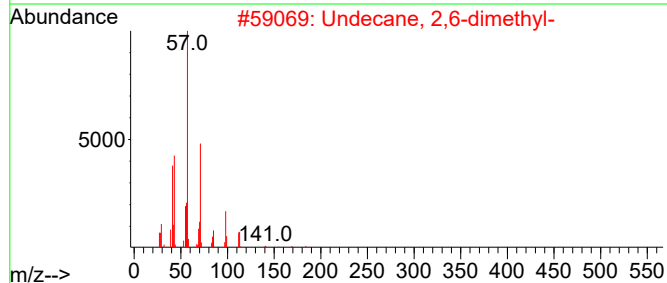
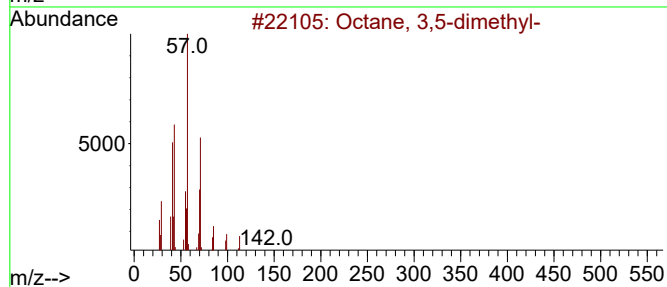
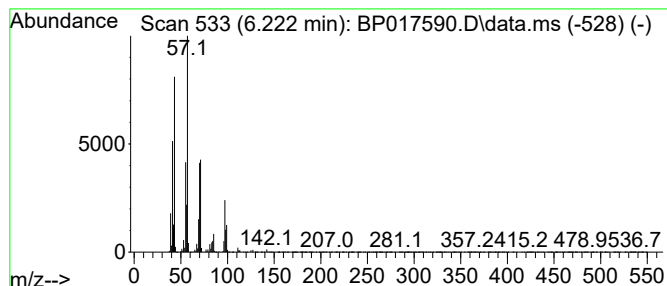
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.222	13.99 ng/ul	3810350	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octane, 3,5-dimethyl-	142	C10H22	015869-93-9	52
2			Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	50
3			1-Butanol, 2-ethyl-	102	C6H14O	000097-95-0	49
4			Nonane, 2-methyl-	142	C10H22	000871-83-0	49
5			1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

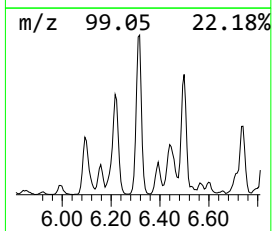
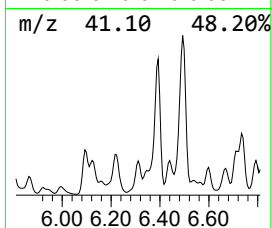
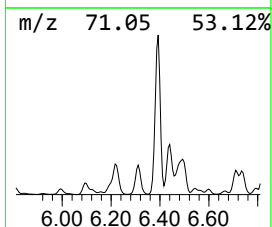
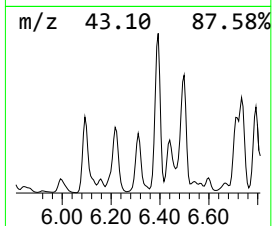
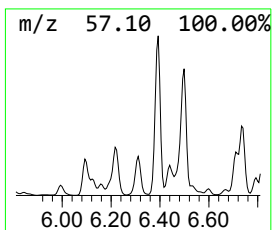
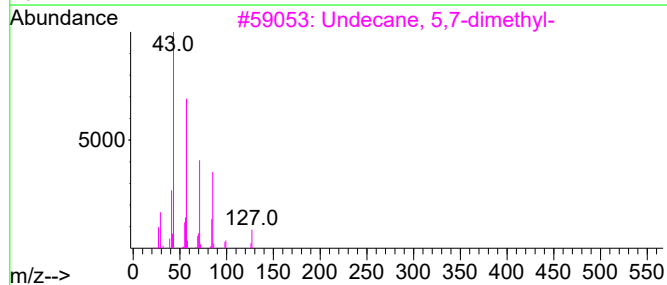
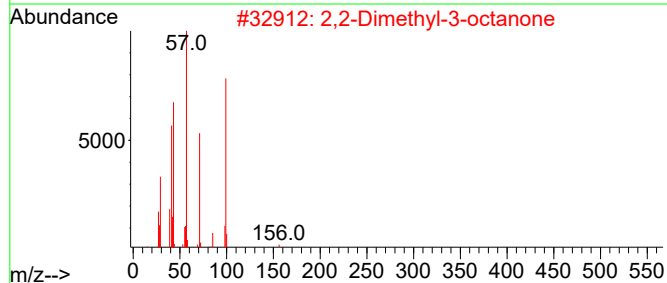
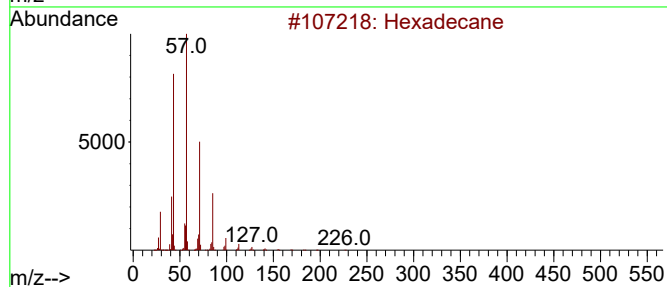
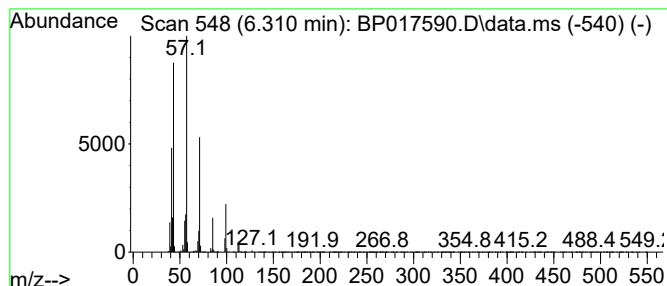
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 (DEL) Alkane: Straight-Chai... Concentration Rank 62

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.310	9.53 ng/ul	2596680	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	64
2	2,2-Dimethyl-3-octanone		156	C10H20O	005340-64-7	59
3	Undecane, 5,7-dimethyl-		184	C13H28	017312-83-3	53
4	3-Heptanone, 5-methyl-		128	C8H16O	000541-85-5	53
5	Undecane, 5-methyl-		170	C12H26	001632-70-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

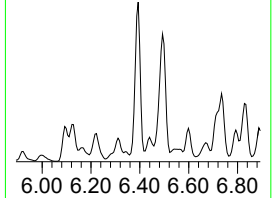
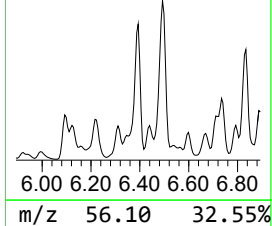
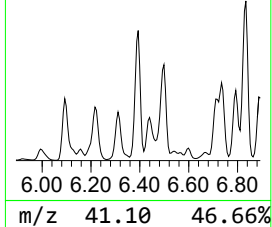
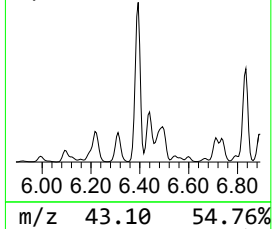
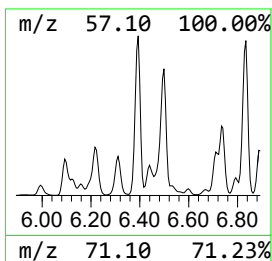
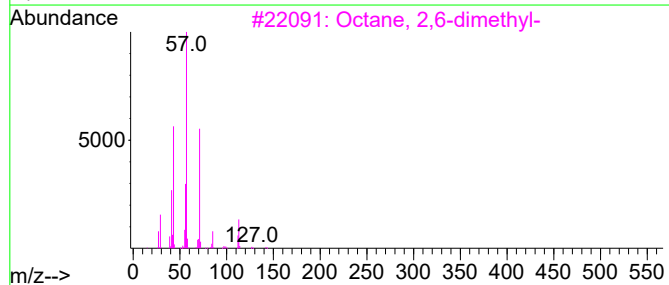
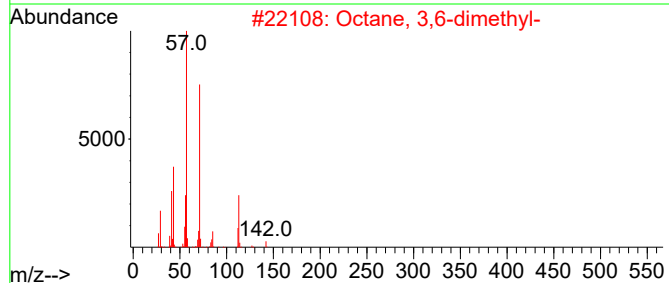
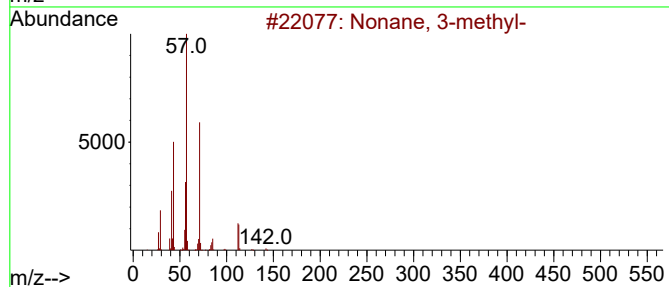
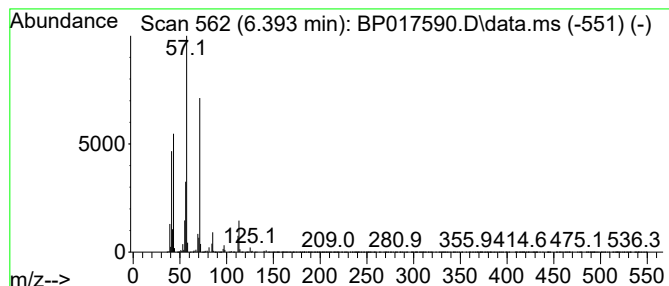
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.393	48.74 ng/ul	13276900	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-		142	C10H22	005911-04-6	91
2	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	91
3	Octane, 2,6-dimethyl-		142	C10H22	002051-30-1	90
4	Octane, 3,6-dimethyl-		142	C10H22	015869-94-0	90
5	Nonane, 3-methyl-		142	C10H22	005911-04-6	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

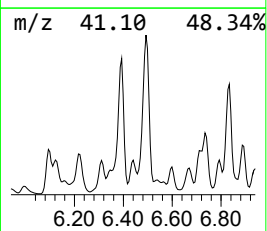
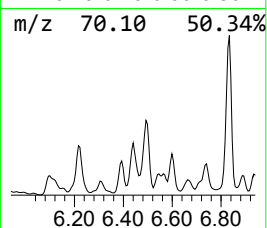
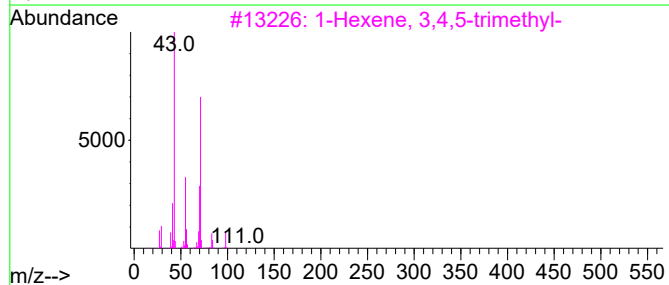
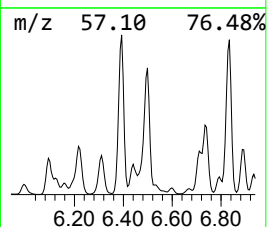
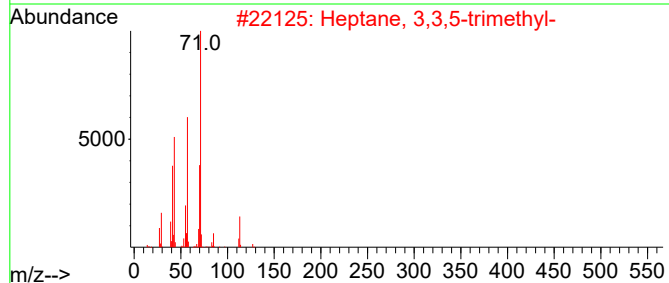
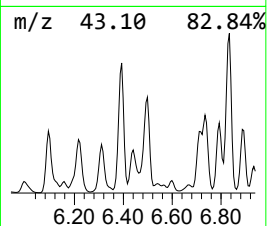
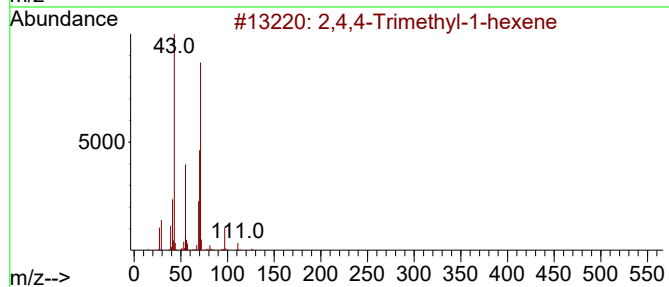
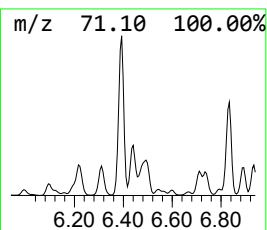
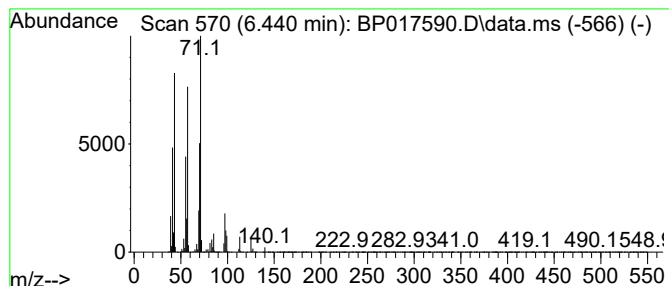
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 63

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.440	9.35 ng/ul	2546280	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,4,4-Trimethyl-1-hexene	126	C9H18	051174-12-0	58
2			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	53
3			1-Hexene, 3,4,5-trimethyl-	126	C9H18	056728-10-0	53
4			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
5			Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

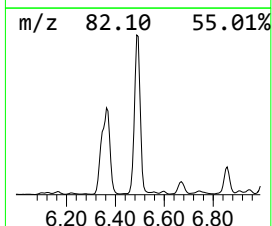
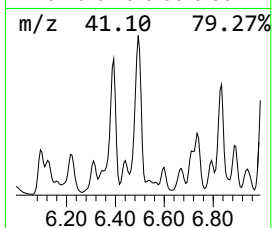
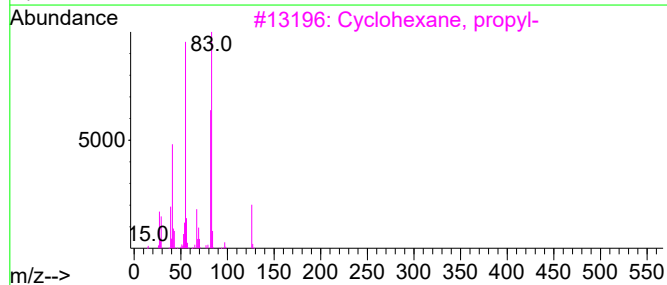
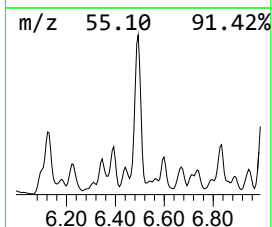
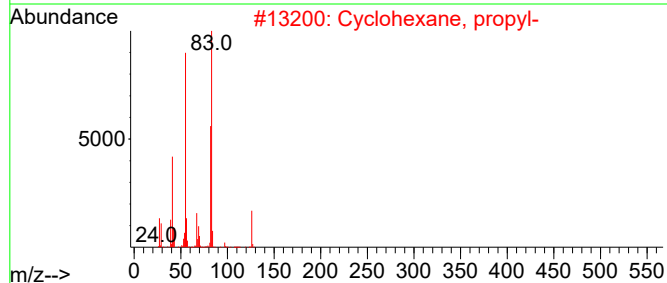
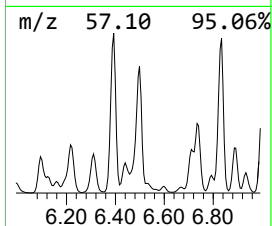
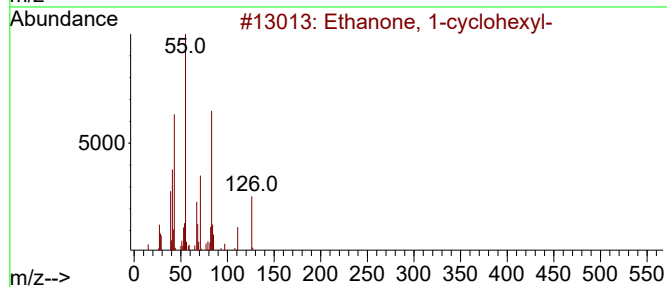
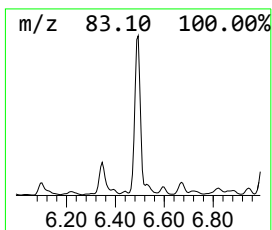
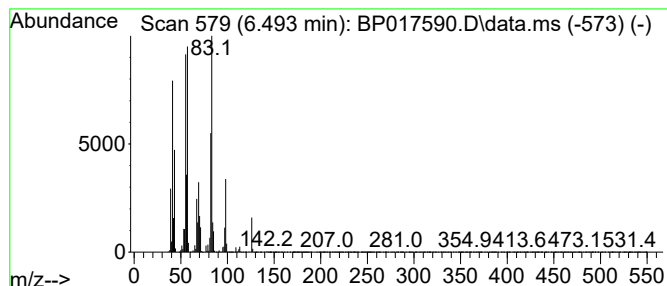
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 13 Ethanone, 1-cyclohexyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.493	53.29 ng/ul	14517000	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanone, 1-cyclohexyl-	126	C8H14O	000823-76-7	53
2			Cyclohexane, propyl-	126	C9H18	001678-92-8	52
3			Cyclohexane, propyl-	126	C9H18	001678-92-8	52
4			Cyclohexane, propyl-	126	C9H18	001678-92-8	52
5			Cyclohexane, methyl-	98	C7H14	000108-87-2	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

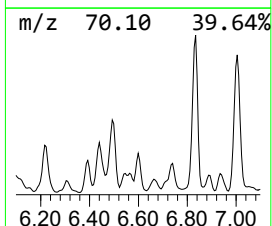
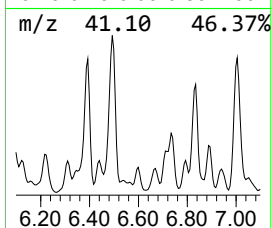
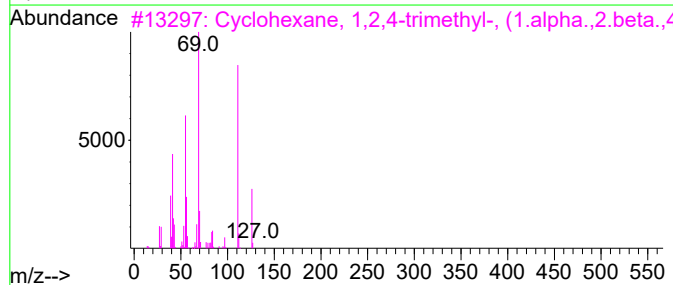
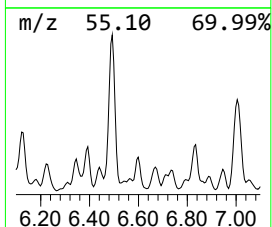
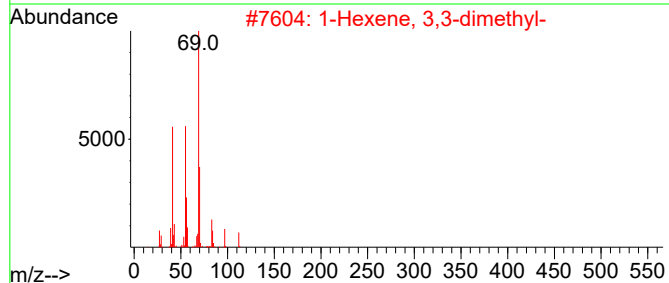
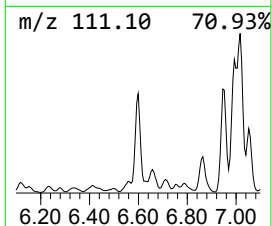
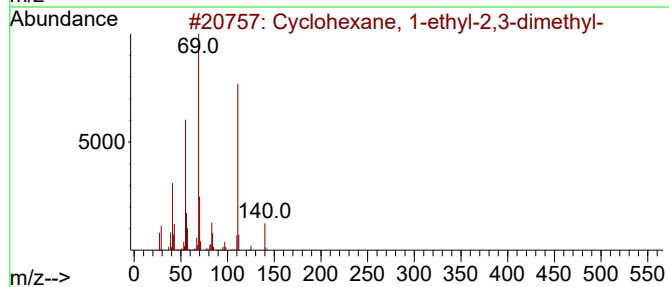
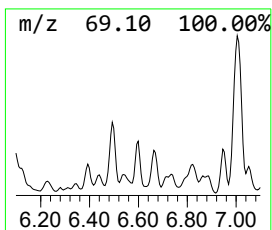
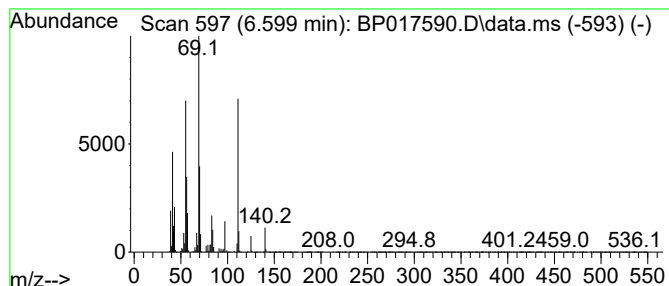
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 14 (DEL) Alkane: Cyclic6.599 Concentration Rank 75

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.599	7.70 ng/ul	2097140	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	72
2			1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	70
3			Cyclohexane, 1,2,4-trimethyl-, (...)	126	C9H18	007667-60-9	53
4			1-Hexene, 3,3,5-trimethyl-	126	C9H18	013427-43-5	50
5			Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

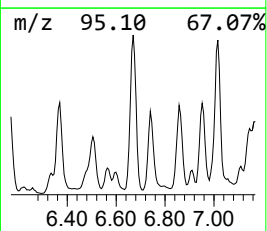
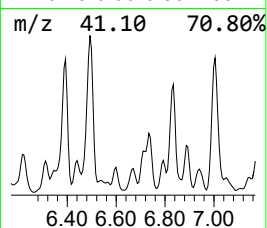
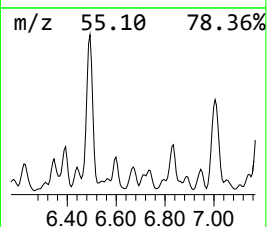
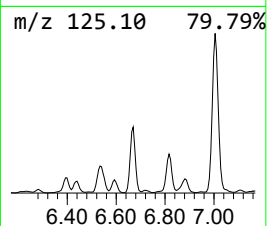
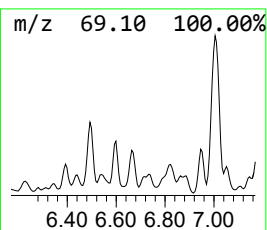
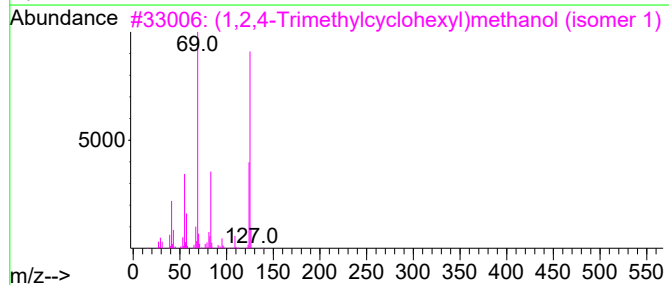
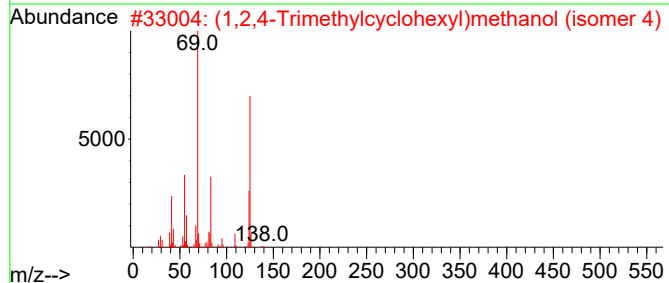
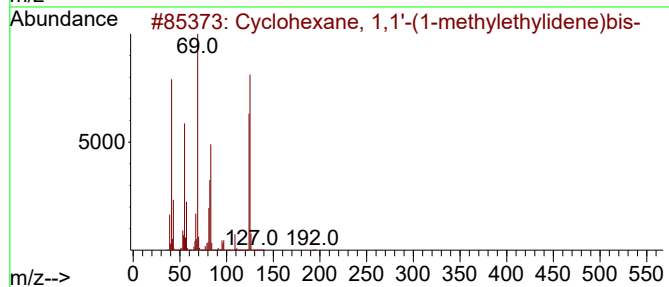
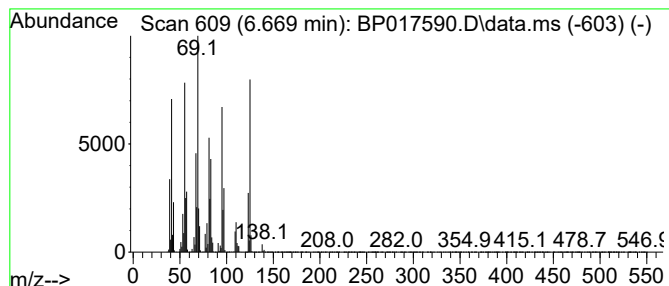
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 15 Cyclohexane, 1,1'-(1-methyl... Concentration Rank 60

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.669	10.42 ng/ul	2837250	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1'-(1-methylethyl...		208	C15H28	054934-90-6	50
2	(1,2,4-Trimethylcyclohexyl)metha...		156	C10H200	1000499-04-1	47
3	(1,2,4-Trimethylcyclohexyl)metha...		156	C10H200	1000499-03-8	47
4	Cyclohexane, 1,1,3,5-tetramethyl...		140	C10H20	050876-32-9	46
5	Cyclohexane, 1,1,3,5-tetramethyl...		140	C10H20	050876-32-9	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

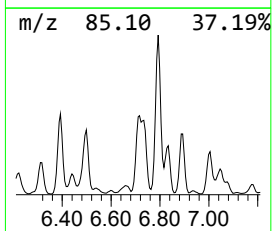
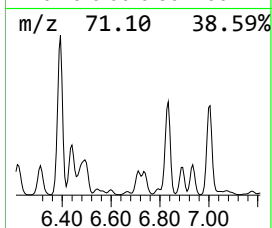
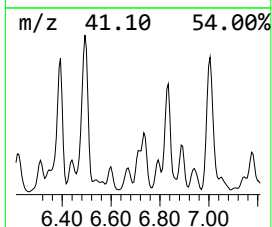
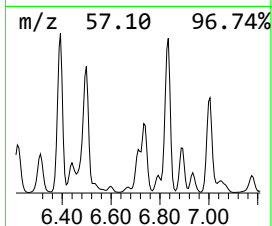
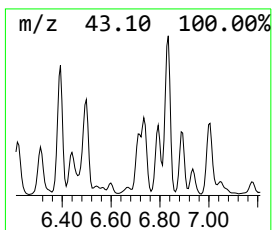
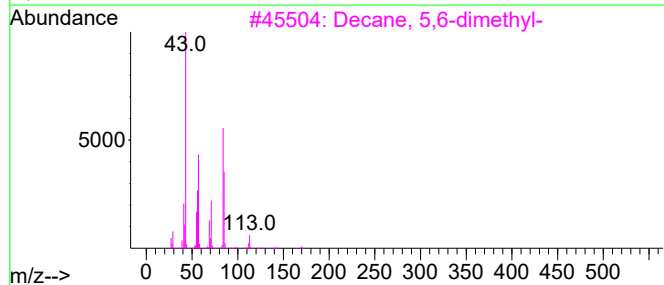
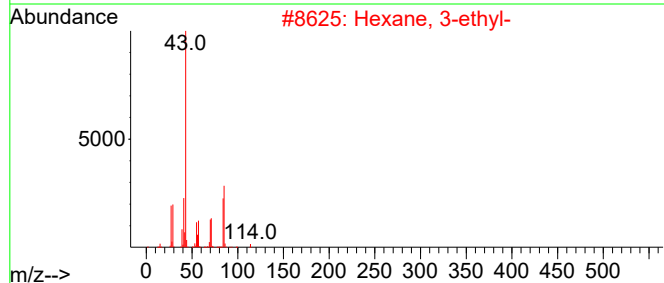
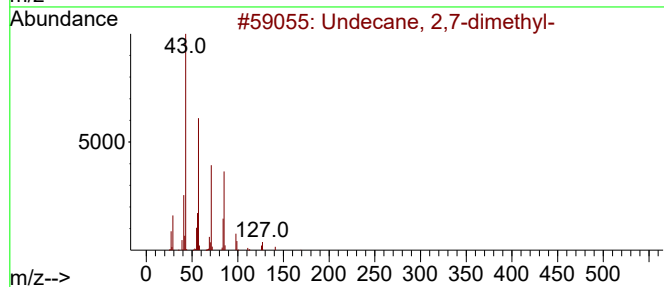
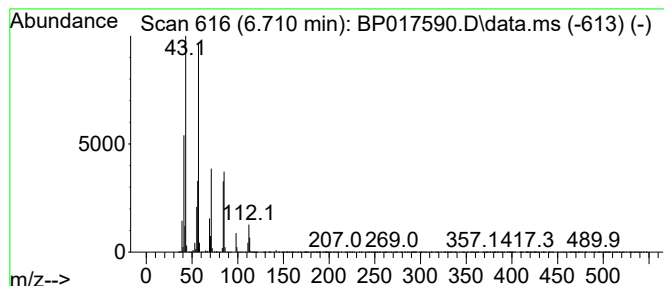
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 72

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.710	8.19 ng/ul	2231510	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 2,7-dimethyl-		184	C13H28	017301-24-5	50
2	Hexane, 3-ethyl-		114	C8H18	000619-99-8	50
3	Decane, 5,6-dimethyl-		170	C12H26	001636-43-7	50
4	1-Iodoundecane		282	C11H23I	004282-44-4	49
5	Nonane, 3-methyl-		142	C10H22	005911-04-6	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

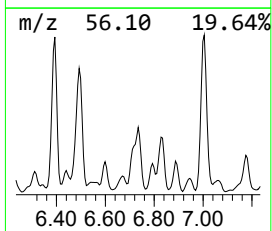
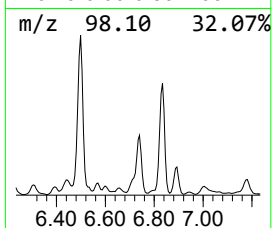
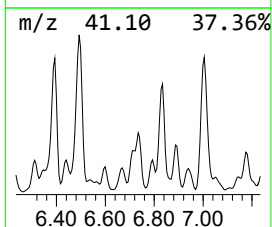
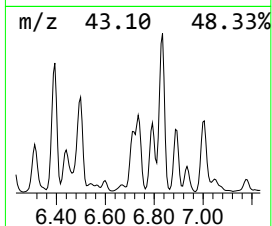
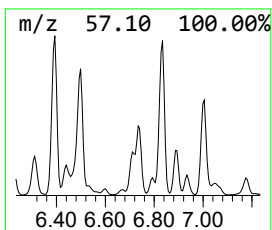
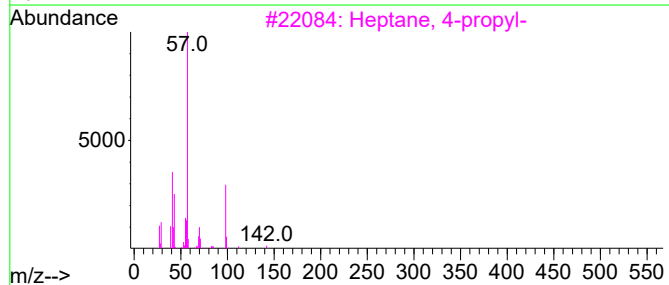
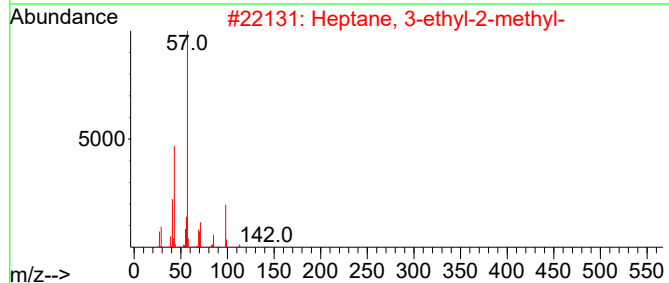
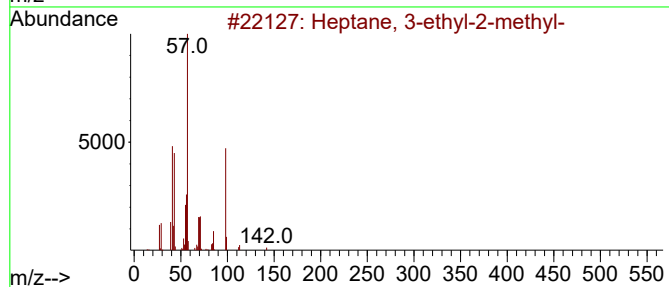
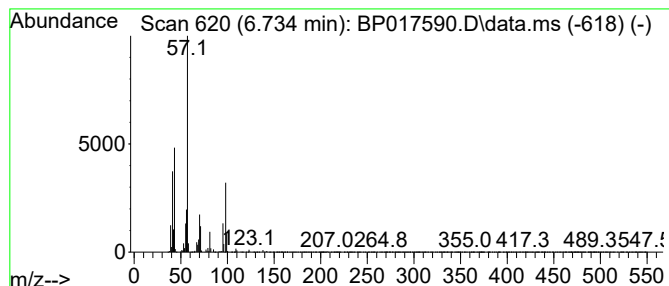
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 17 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.734	13.18 ng/ul	3588900	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	59
2	Heptane, 3-ethyl-2-methyl-	142	C10H22	014676-29-0	53
3	Heptane, 4-propyl-	142	C10H22	003178-29-8	50
4	Octane, 3-methyl-	128	C9H20	002216-33-3	50
5	1,3,2-Dioxaborolane, 4,4-dimethy...	142	C6H11BO3	1000063-89-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

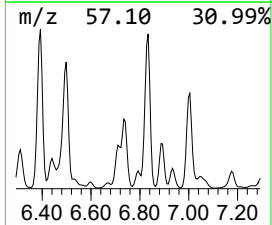
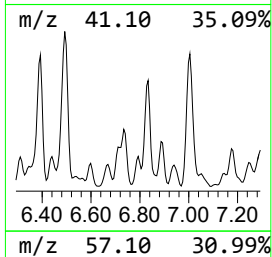
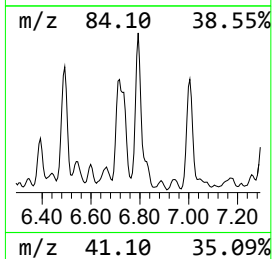
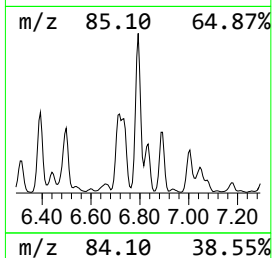
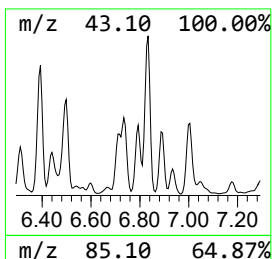
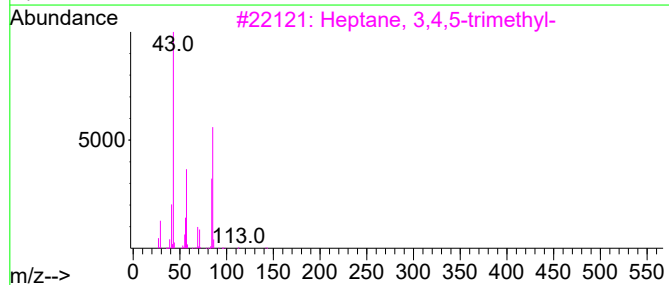
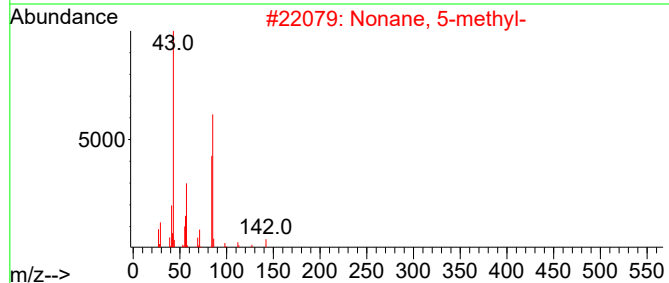
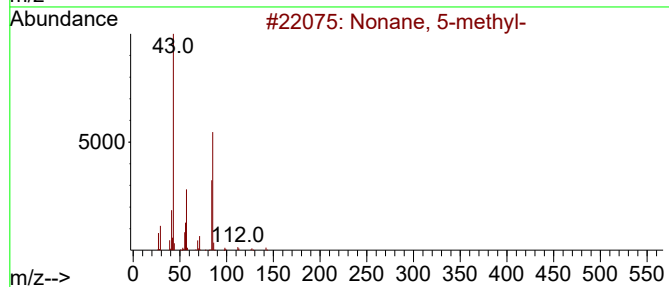
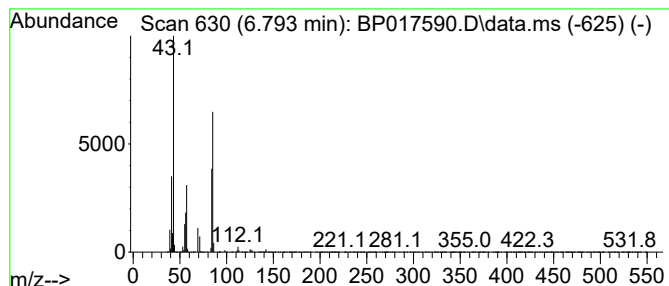
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 74

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.793	7.77 ng/ul	2115480	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 5-methyl-	142	C10H22	015869-85-9	91
2			Nonane, 5-methyl-	142	C10H22	015869-85-9	91
3			Heptane, 3,4,5-trimethyl-	142	C10H22	020278-89-1	90
4			Pentane, 3-ethyl-2,4-dimethyl-	128	C9H20	001068-87-7	72
5			Hexane, 2,3,5-trimethyl-	128	C9H20	001069-53-0	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

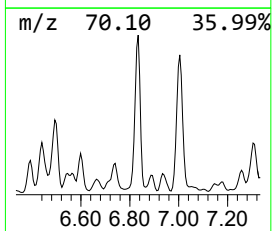
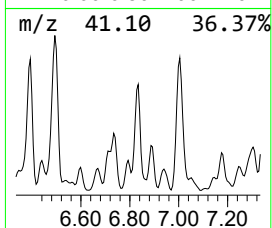
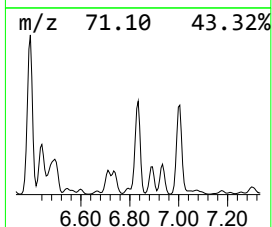
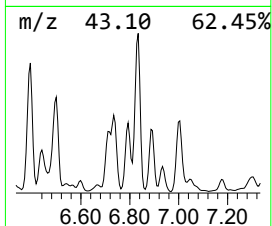
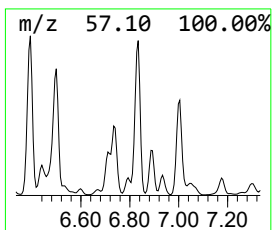
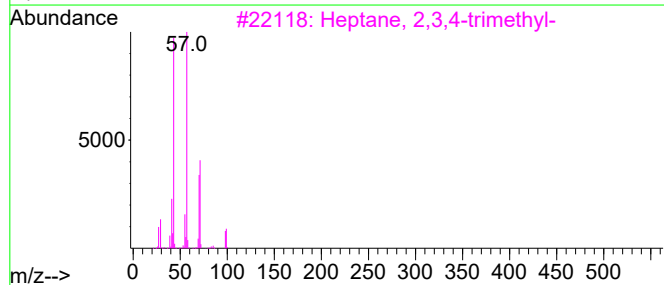
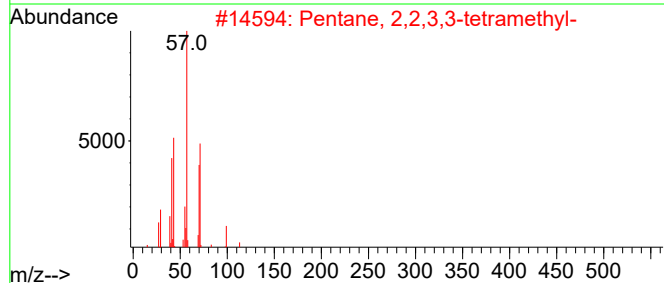
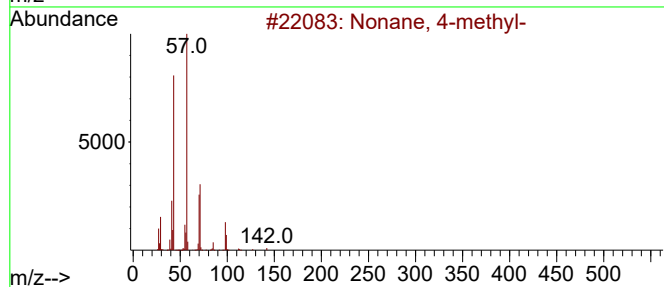
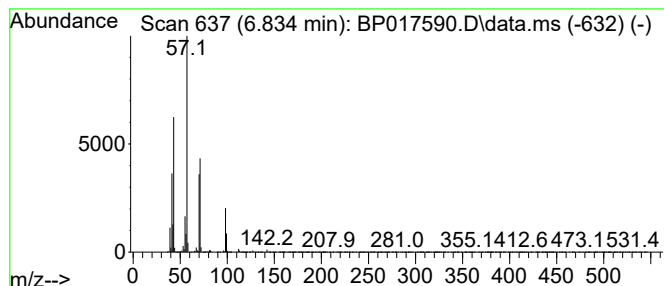
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD			R.T.
6.834	30.79 ng/ul	8388620	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-		142	C10H22	017301-94-9	87
2	Pentane, 2,2,3,3-tetramethyl-		128	C9H20	007154-79-2	59
3	Heptane, 2,3,4-trimethyl-		142	C10H22	052896-95-4	59
4	Undecane, 4,5-dimethyl-		184	C13H28	017312-79-7	59
5	Hexane, 3-methyl-		100	C7H16	000589-34-4	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

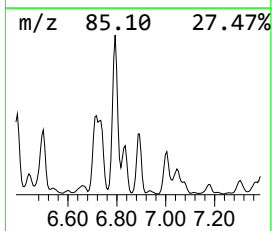
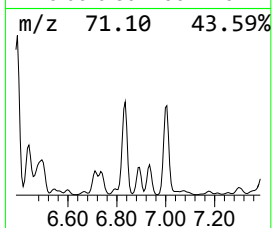
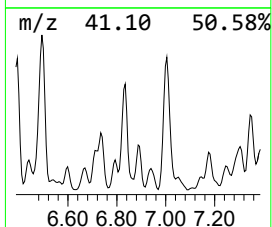
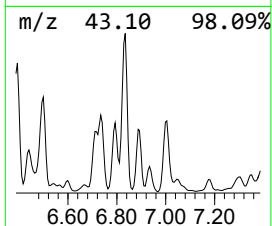
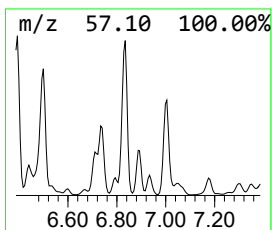
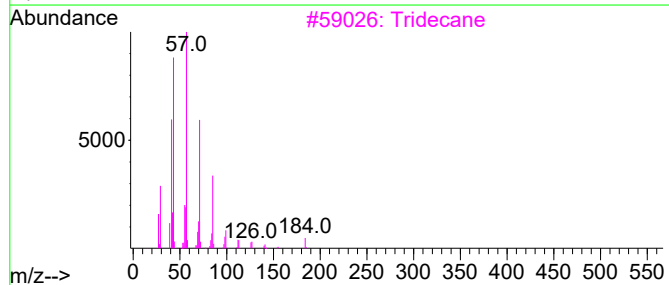
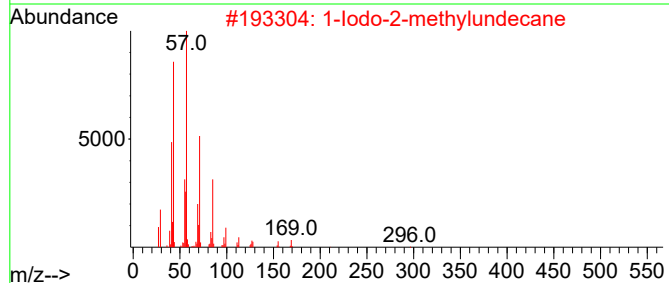
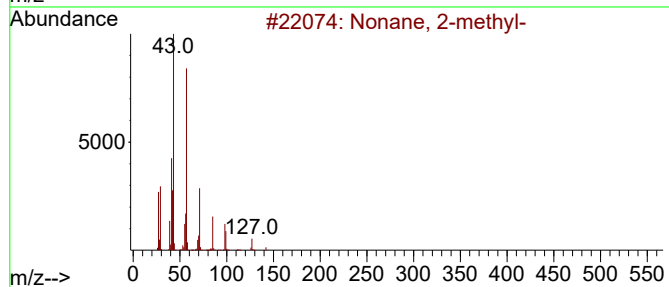
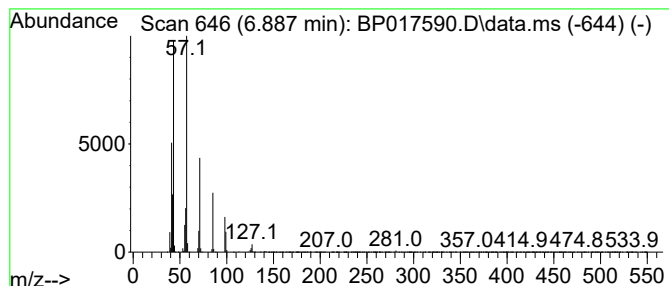
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 64

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.887	9.04 ng/ul	2462560	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Nonane, 2-methyl-	142	C10H22	000871-83-0	83
2		1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	78
3		Tridecane	184	C13H28	000629-50-5	72
4		Dodecane	170	C12H26	000112-40-3	72
5		Nonane, 2-methyl-	142	C10H22	000871-83-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

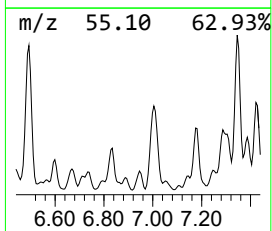
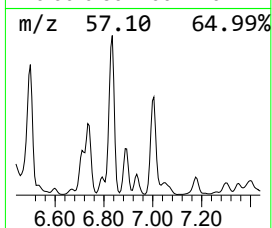
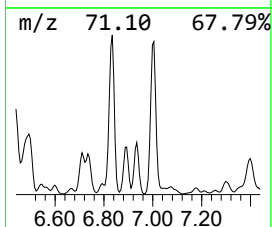
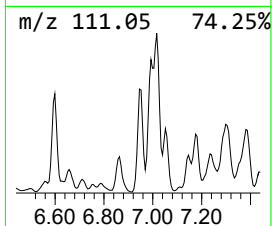
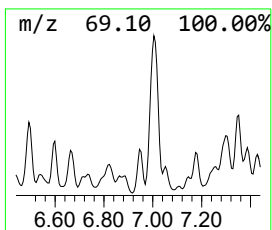
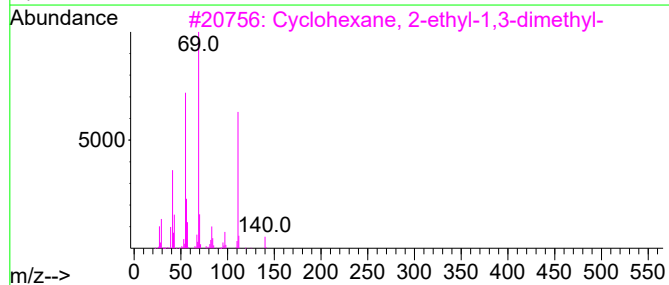
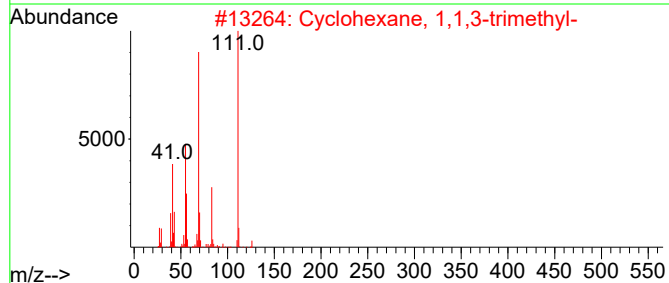
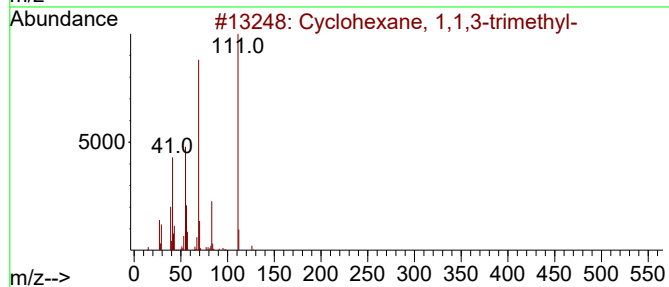
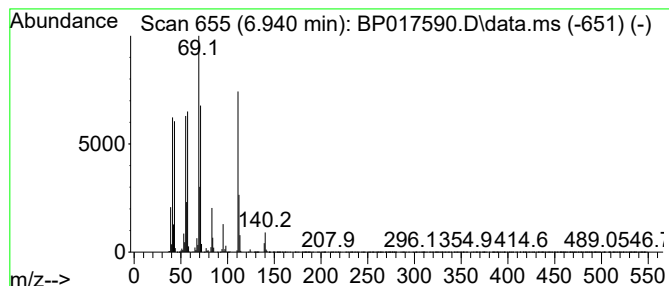
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 21 (DEL) Alkane: Cyclic6.940 Concentration Rank 70

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.940	8.23 ng/ul	2243060	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	53
2			Cyclohexane, 1,1,3-trimethyl-	126	C9H18	003073-66-3	53
3			Cyclohexane, 2-ethyl-1,3-dimethyl-	140	C10H20	007045-67-2	52
4			Cyclohexane, 1-ethyl-2,3-dimethyl-	140	C10H20	007058-05-1	50
5			Cyclohexane, 1,3,5-trimethyl-, (...)	126	C9H18	001795-26-2	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

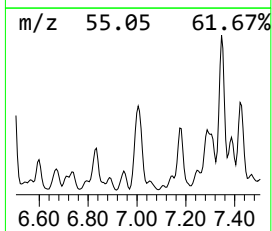
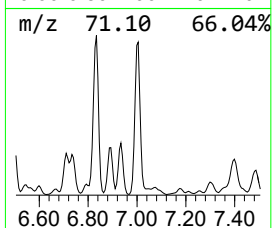
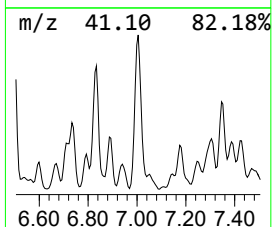
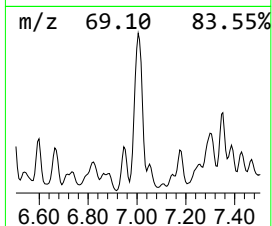
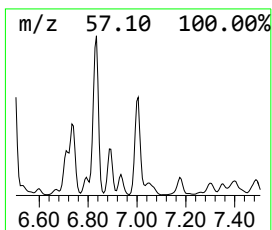
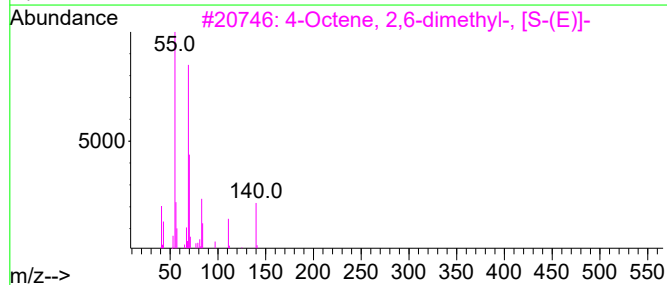
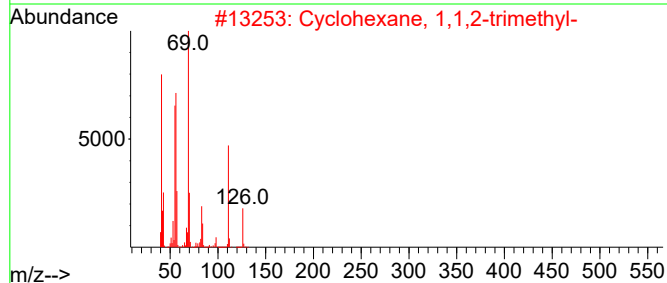
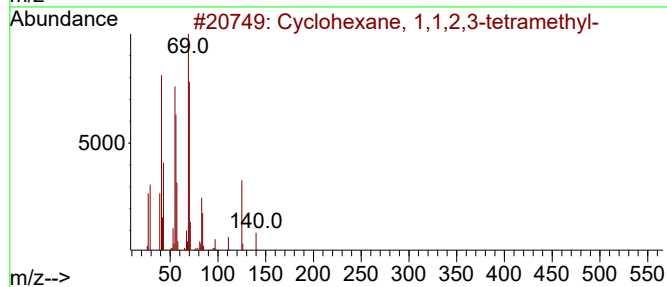
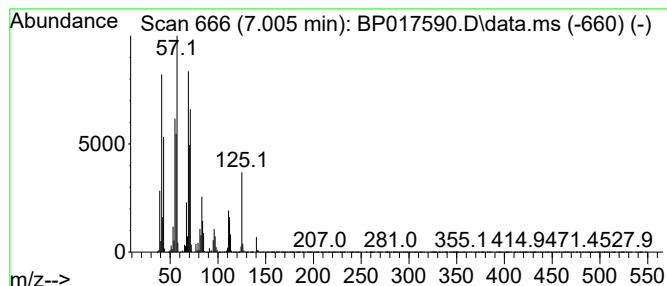
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 22 (DEL) Alkane: Cyclic7.005 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.005	52.33 ng/ul	14256000	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1,1,2,3-tetramethyl-	140	C10H20	006783-92-2	70
2		Cyclohexane, 1,1,2-trimethyl-	126	C9H18	007094-26-0	60
3		4-Octene, 2,6-dimethyl-, [S-(E)]-	140	C10H20	062960-76-3	53
4		Isodecyl methacrylate	226	C14H26O2	029964-84-9	47
5		2-Octene, 2,6-dimethyl-	140	C10H20	004057-42-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

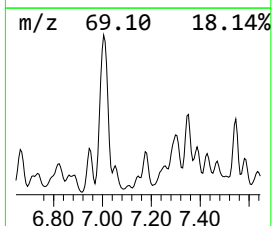
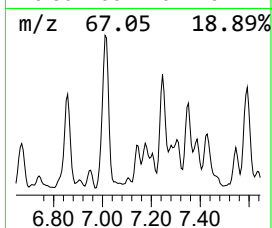
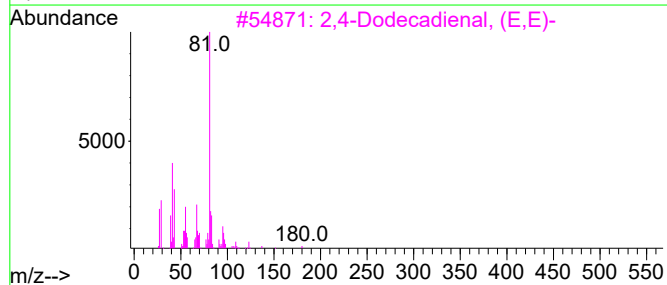
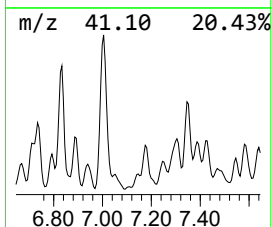
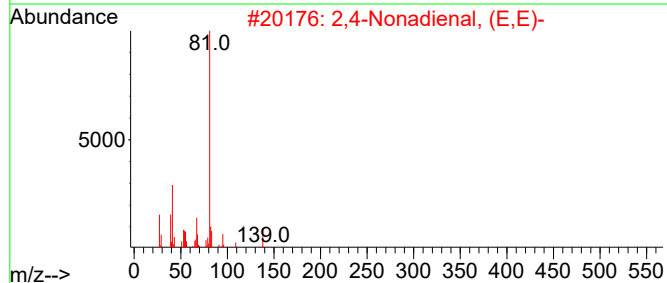
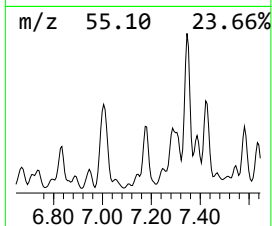
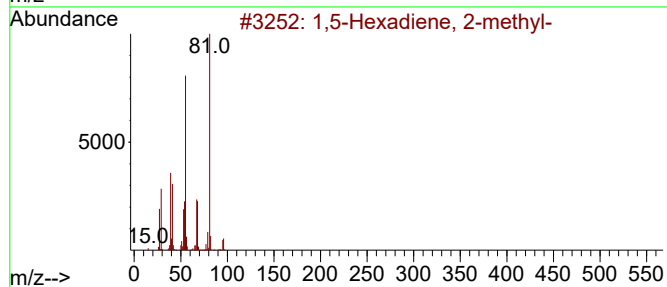
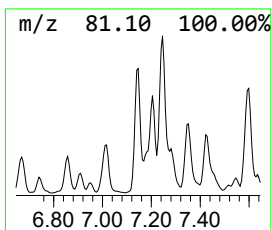
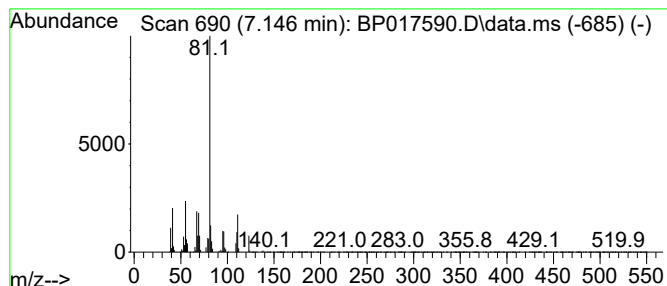
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 23 1,5-Hexadiene, 2-methyl- Concentration Rank 78

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.146	5.97 ng/ul	1626940	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,5-Hexadiene, 2-methyl-	96	C7H12	004049-81-4	53
2			2,4-Nonadienal, (E,E)-	138	C9H14O	005910-87-2	53
3			2,4-Dodecadienal, (E,E)-	180	C12H20O	021662-16-8	45
4			2,4-Dodecadienal	180	C12H20O	013162-47-5	45
5			2,4-Heptadienal, (E,E)-	110	C7H10O	004313-03-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

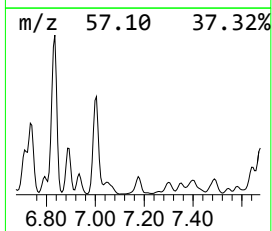
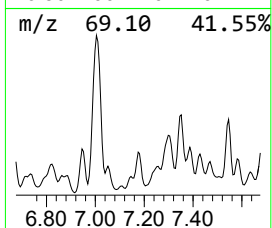
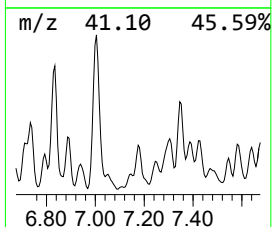
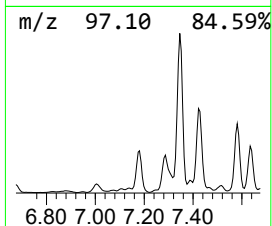
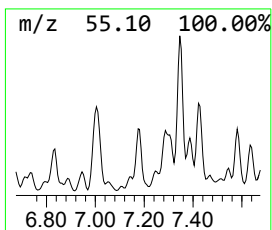
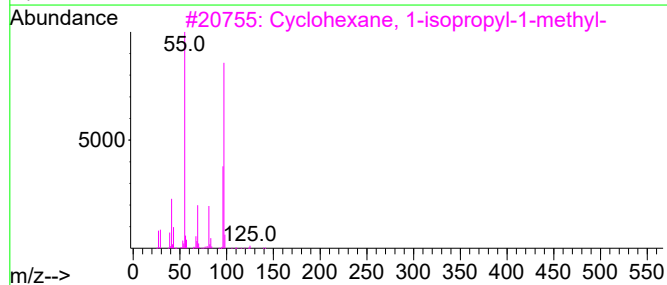
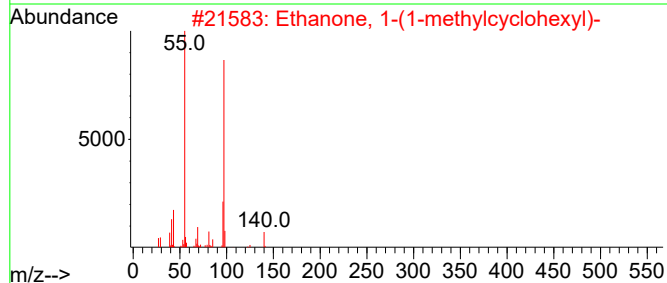
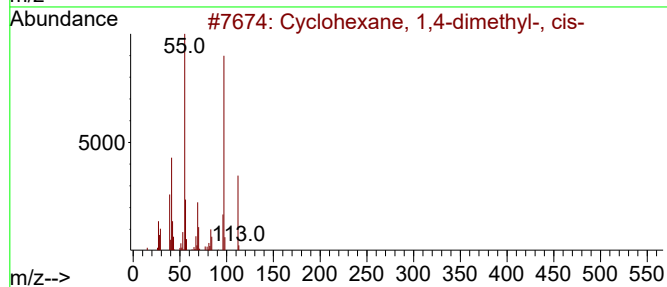
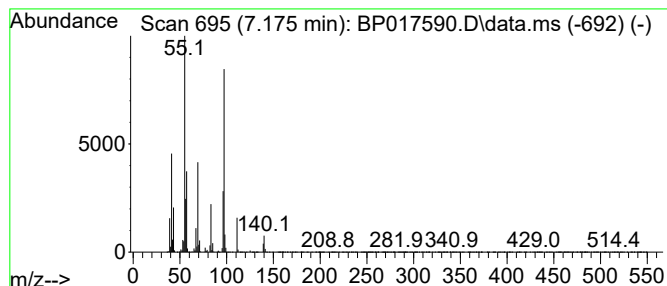
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 24 (DEL) Alkane: Cyclic7.175 Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.175	15.03 ng/ul	4093370	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-, cis-	112	C8H16	000624-29-3	59
2			Ethanone, 1-(1-methylcyclohexyl)-	140	C9H16O	002890-62-2	58
3			Cyclohexane, 1-isopropyl-1-methyl-	140	C10H20	016580-26-0	53
4			Cyclohexane, 1-ethyl-2-methyl-	126	C9H18	003728-54-9	53
5			Oxalic acid, cyclohexylmethyl et...	214	C11H18O4	1000309-68-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

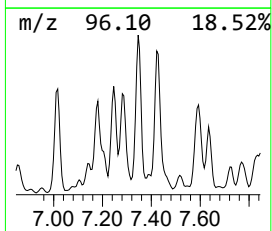
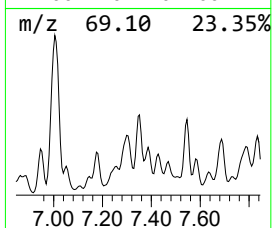
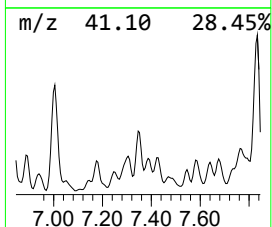
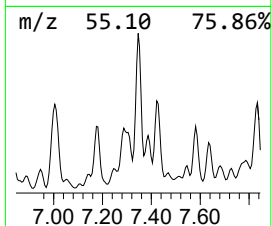
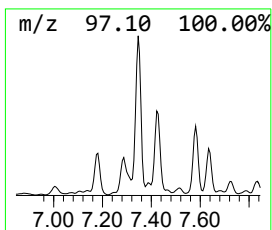
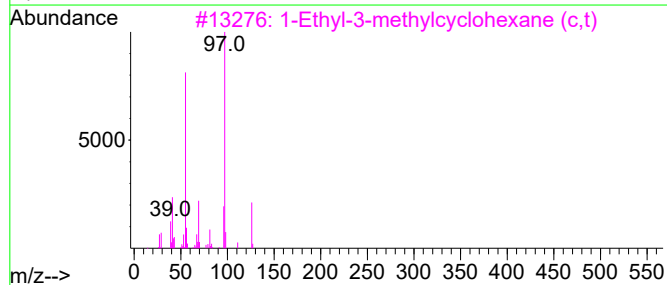
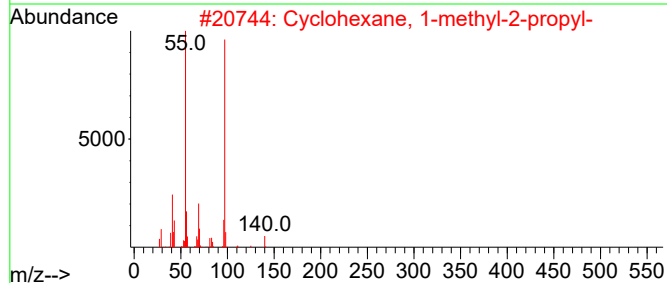
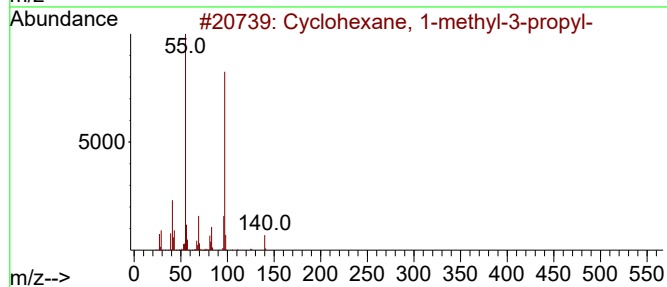
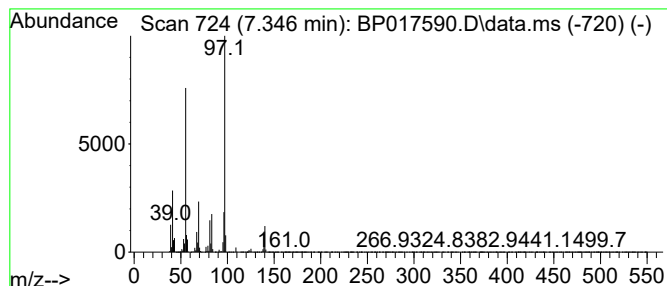
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 25 (DEL) Alkane: Cyclic7.346 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.346	30.29 ng/ul	8249760	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	74
2			Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	72
3			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	64
4			Cyclohexane, 1-methyl-4-(1-methyl-...	140	C10H20	001678-82-6	64
5			Cyclohexane, 1-bromo-4-methyl-	176	C7H13Br	006294-40-2	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

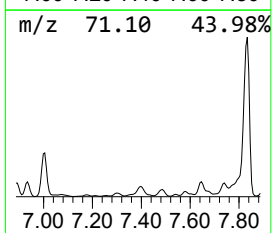
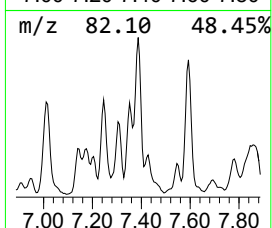
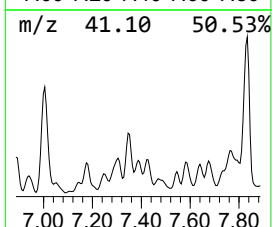
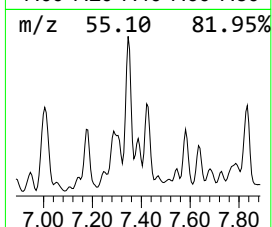
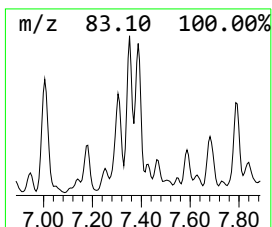
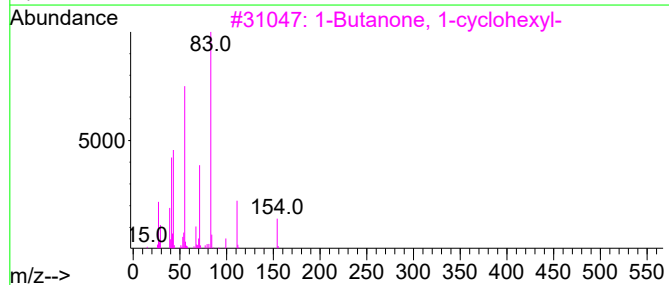
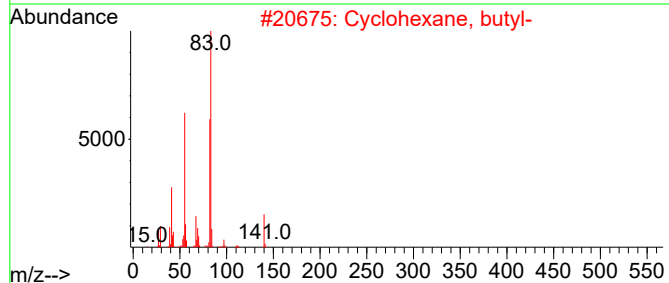
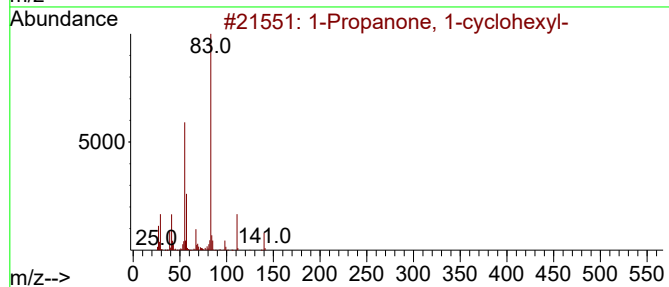
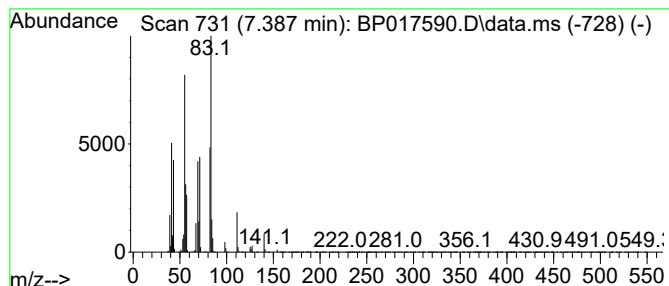
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 26 1-Propanone, 1-cyclohexyl- Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.387	12.14 ng/ul	3305670	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Propanone, 1-cyclohexyl-	140	C9H16O	001123-86-0	60
2	Cyclohexane, butyl-	140	C10H20	001678-93-9	52
3	1-Butanone, 1-cyclohexyl-	154	C10H18O	001462-27-7	50
4	Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	50
5	Cyclohexane, propyl-	126	C9H18	001678-92-8	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

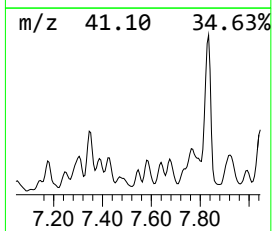
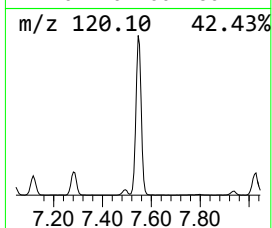
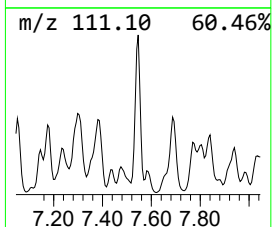
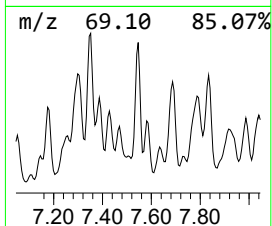
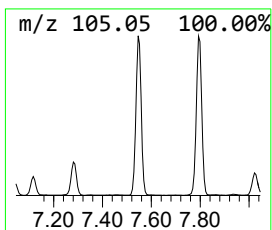
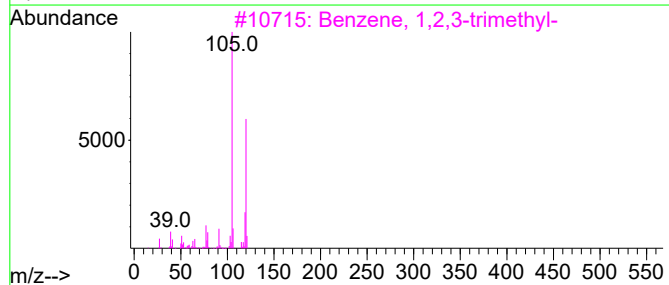
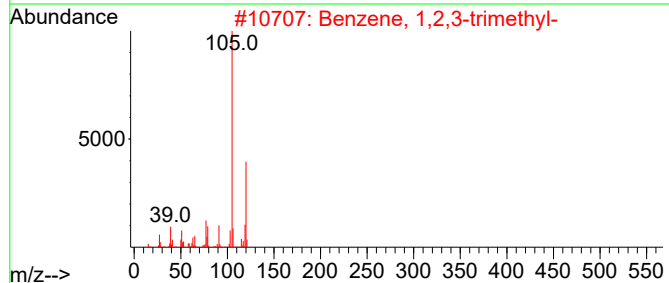
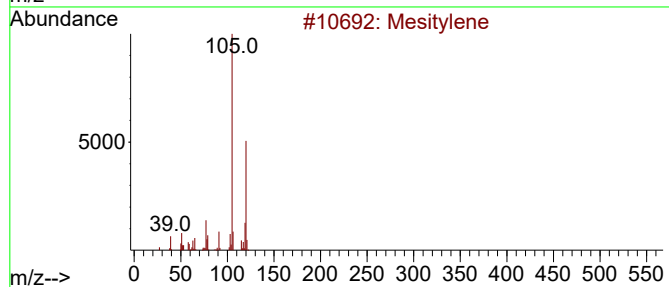
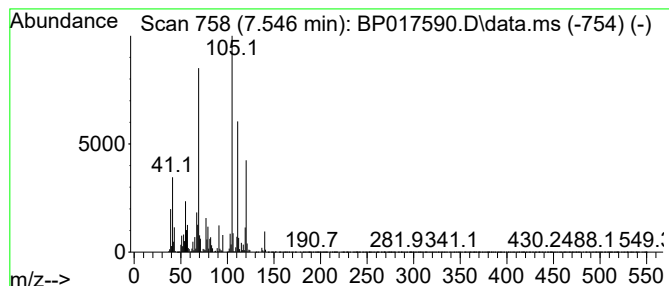
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 27 Mesitylene Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.546	10.43 ng/ul	2841390	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	78
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	74
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	74
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	70
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

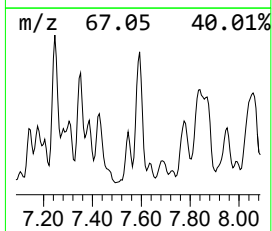
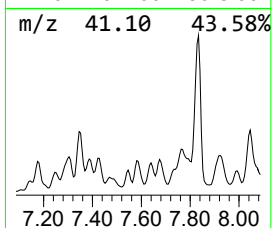
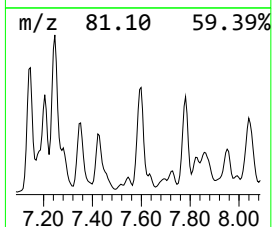
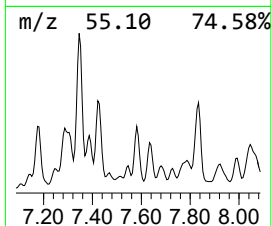
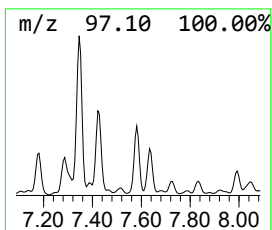
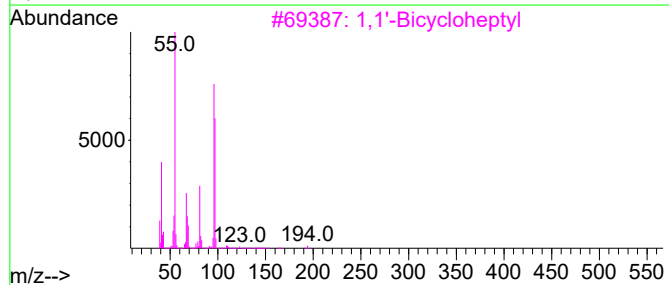
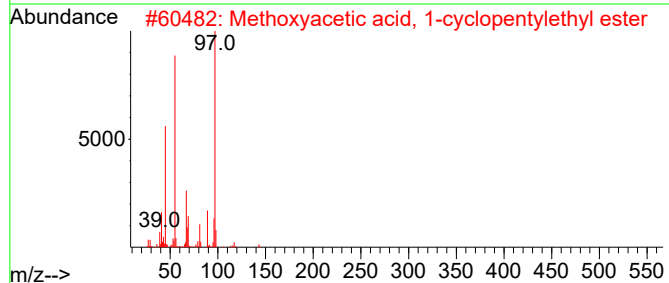
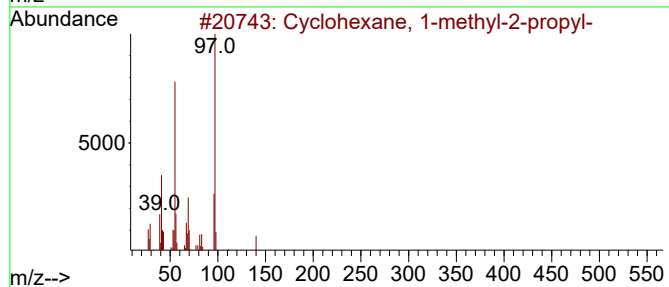
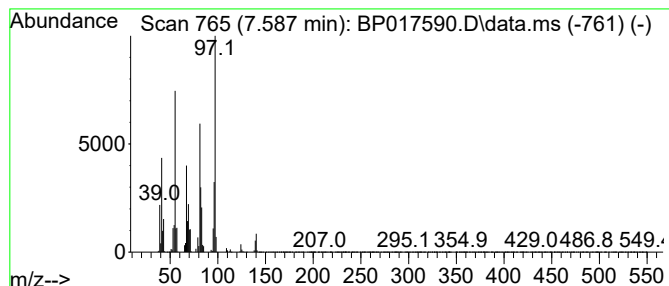
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 28 (DEL) Alkane: Cyclic7.587 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.587	11.94 ng/ul	3251950	1,4-Dichlorobenzene-d4			7.940
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 1-methyl-2-propyl-	140	C10H20	004291-79-6	53
2		Methoxyacetic acid, 1-cyclopenty...	186	C10H18O3	1000282-69-5	50
3		1,1'-Bicycloheptyl	194	C14H26	023183-11-1	47
4		Cyclohexane, 1-methyl-3-(1-methy...	140	C10H20	016580-24-8	47
5		Sulfurous acid, cyclohexylmethyl...	262	C13H26O3S	1000309-21-5	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

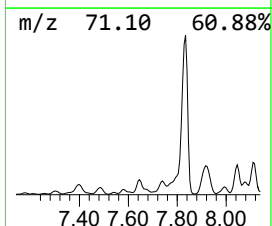
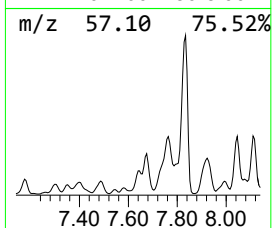
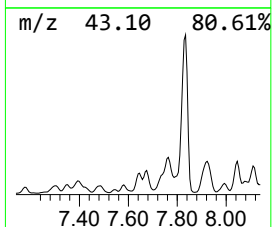
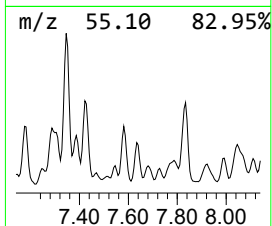
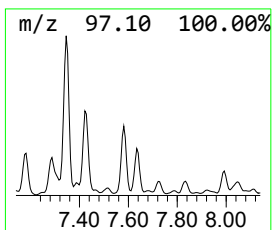
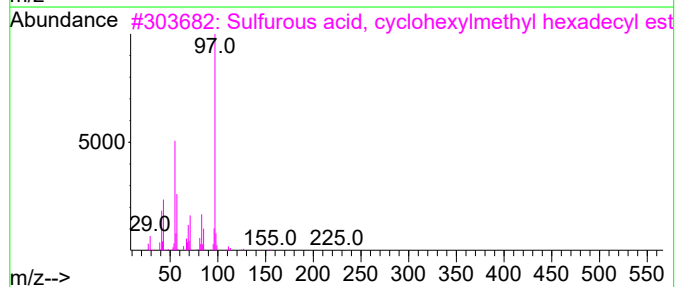
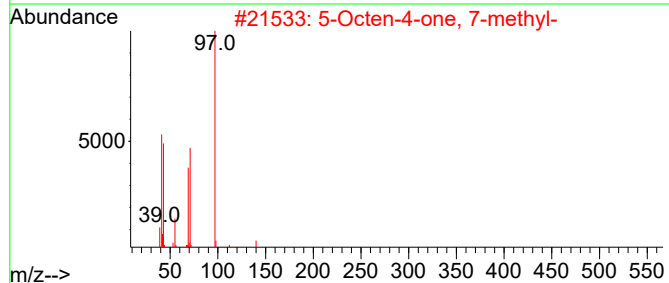
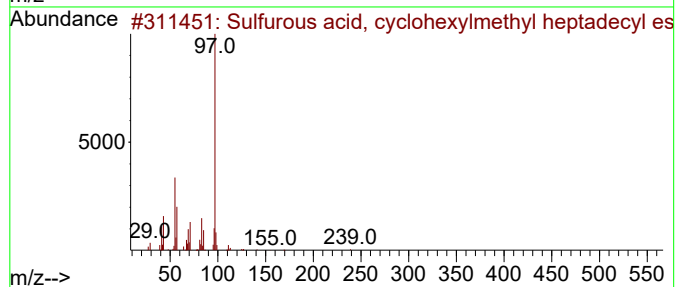
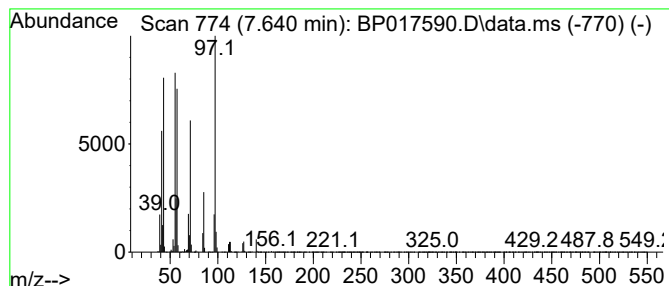
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 29 Sulfurous acid, cyclohexylm... Concentration Rank 68

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.640	8.47 ng/ul	2307220	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Sulfurous acid, cyclohexylmethyl...	416	C24H48O3S	1000309-22-5	53
2			5-Octen-4-one, 7-methyl-	140	C9H16O	032064-78-1	53
3			Sulfurous acid, cyclohexylmethyl...	402	C23H46O3S	1000309-22-4	50
4			Sulfurous acid, cyclohexylmethyl...	374	C21H42O3S	1000309-22-2	50
5			Cyclohexane, 1-methyl-3-propyl-	140	C10H20	004291-80-9	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

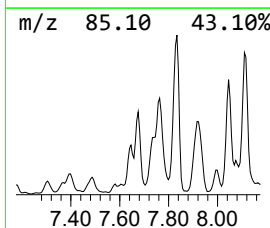
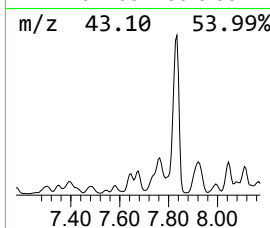
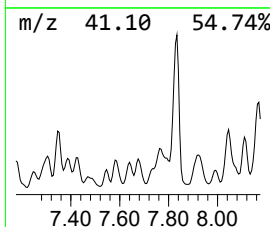
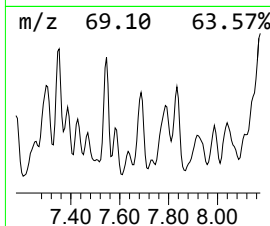
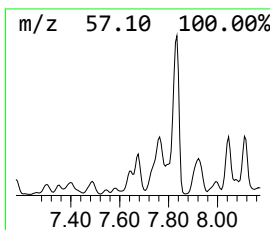
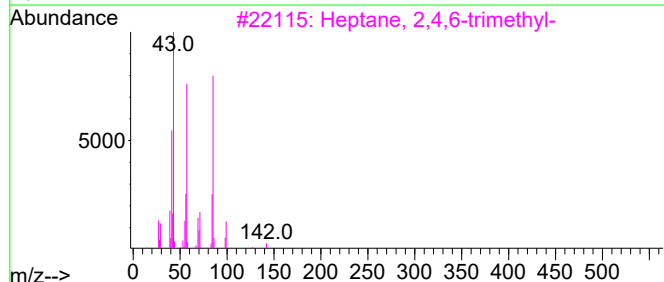
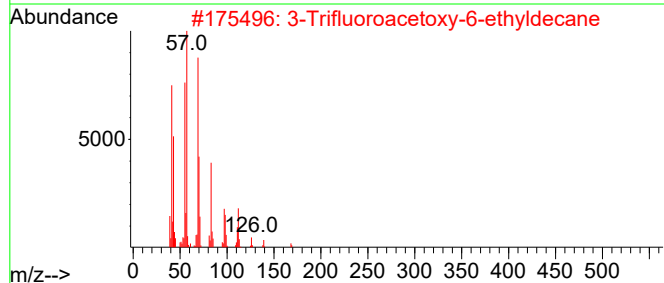
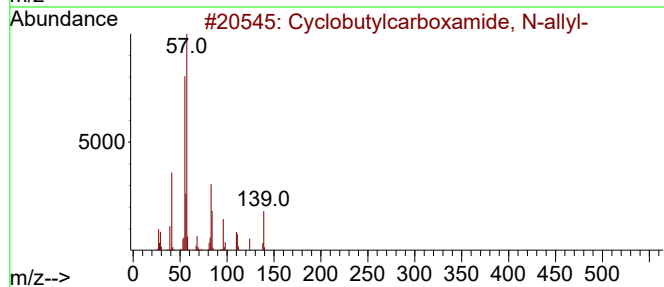
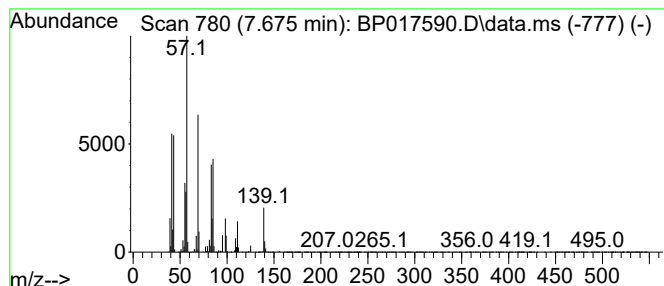
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 30 Cyclobutylcarboxamide, N-al... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.675	9.66 ng/ul	2632230	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclobutylcarboxamide, N-allyl-	139	C8H13NO	1000340-36-9	50
2			3-Trifluoroacetoxy-6-ethyldecane	282	C14H25F3O2	116436-59-0	43
3			Heptane, 2,4,6-trimethyl-	142	C10H22	002613-61-8	38
4			Decane	142	C10H22	000124-18-5	38
5			Carbonic acid, isobutyl isohexyl...	202	C11H22O3	1000314-60-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

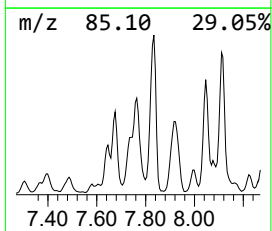
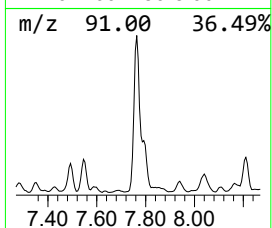
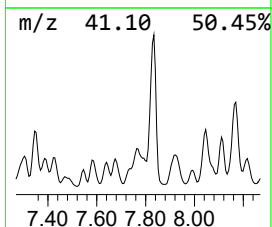
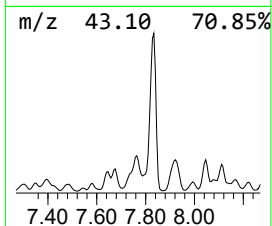
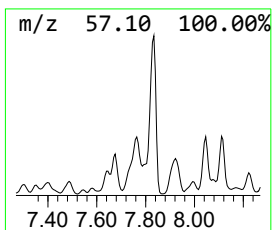
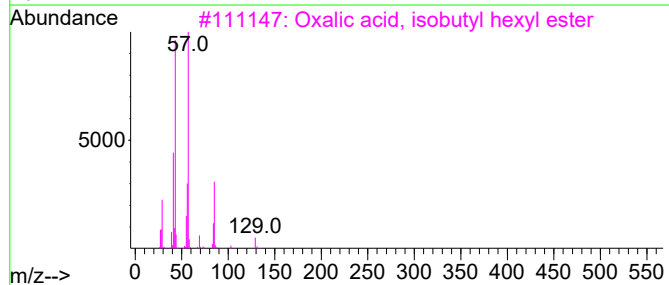
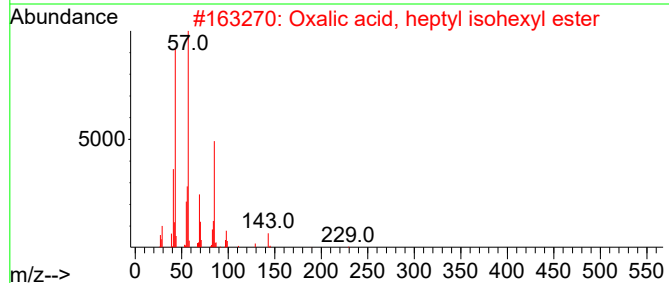
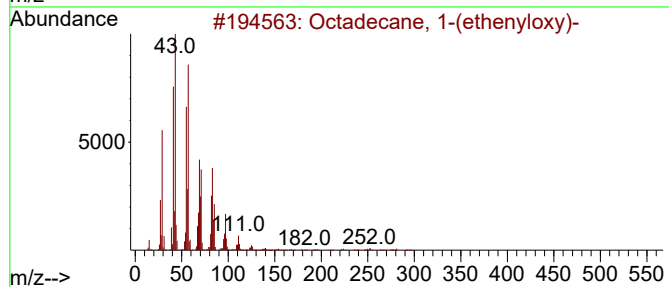
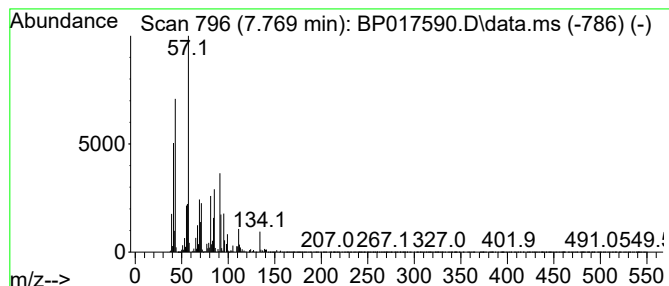
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 31 unknown-01 Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.769	25.76 ng/ul	7016990	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecane, 1-(ethenyloxy)-	296	C20H40O	000930-02-9	43
2			Oxalic acid, heptyl isohexyl ester	272	C15H28O4	1000309-33-1	38
3			Oxalic acid, isobutyl hexyl ester	230	C12H22O4	1000309-37-1	38
4			Octane	114	C8H18	000111-65-9	38
5			Octane	114	C8H18	000111-65-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

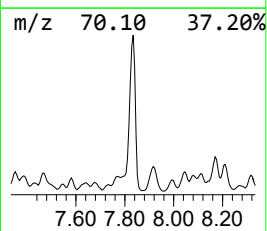
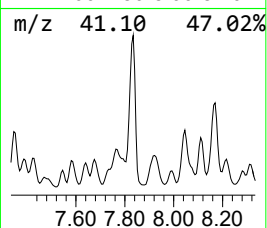
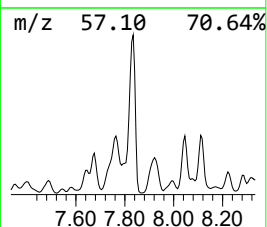
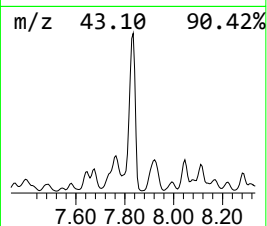
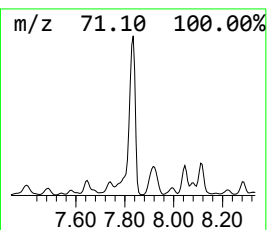
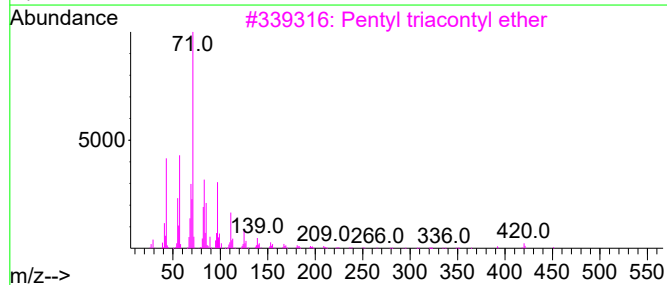
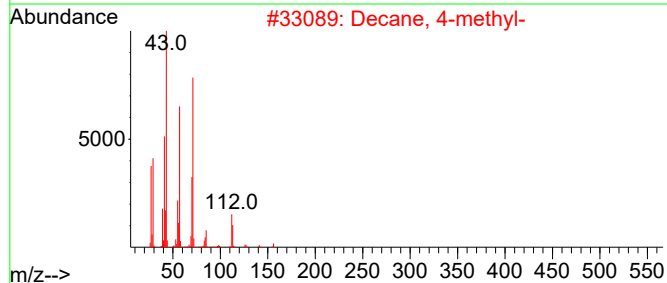
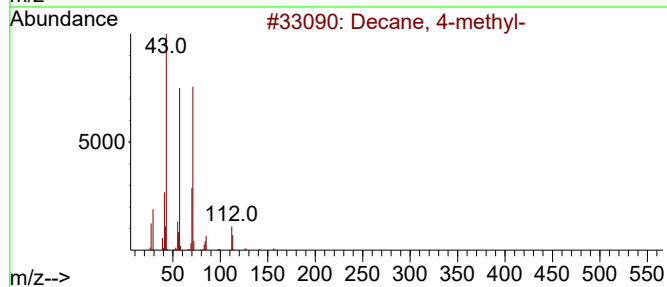
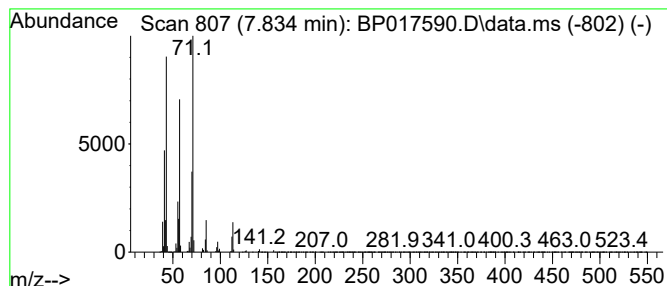
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.834	64.85 ng/ul	17666500	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 4-methyl-	156	C11H24	002847-72-5	90
2	Decane, 4-methyl-	156	C11H24	002847-72-5	78
3	Pentyl triacontyl ether	509	C35H72O	1000406-42-5	72
4	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
5	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

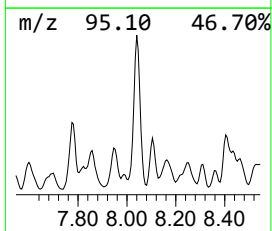
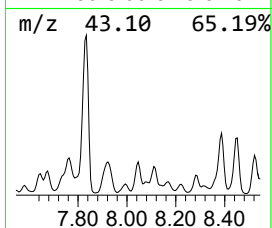
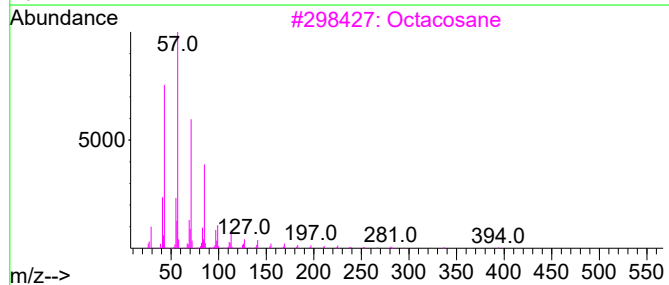
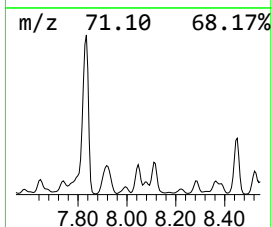
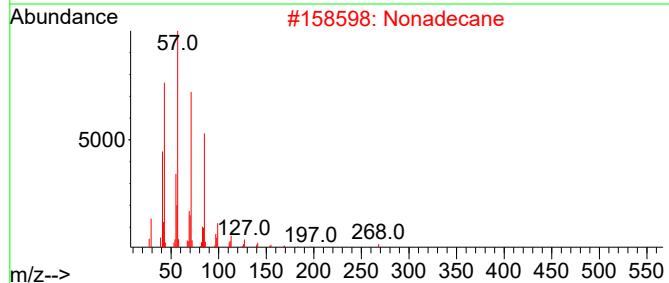
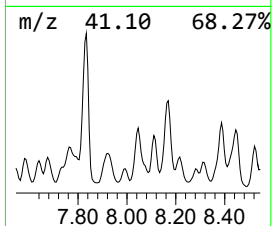
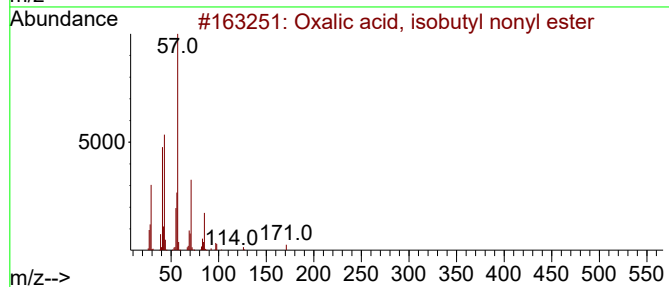
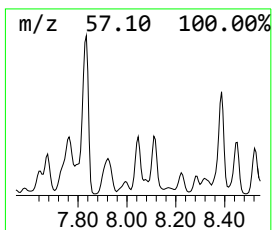
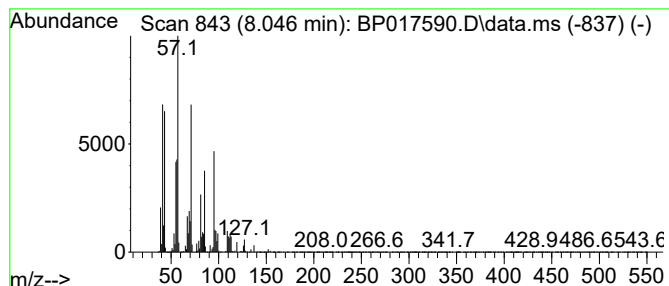
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 33 unknown-02 Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.046	29.66 ng/ul	8080700	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, isobutyl nonyl ester	272	C15H28O4	1010309-37-4	46
2			Nonadecane	268	C19H40	000629-92-5	45
3			Octacosane	394	C28H58	000630-02-4	43
4			Tetracosane	338	C24H50	000646-31-1	43
5			Nonane, 5-(1-methylpropyl)-	184	C13H28	062185-54-0	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

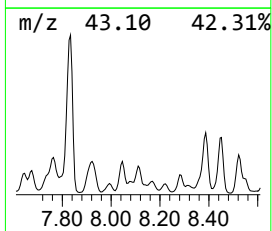
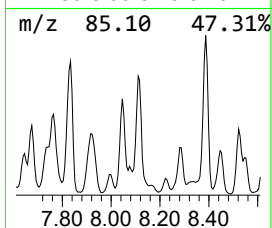
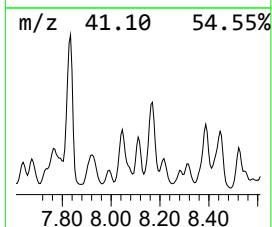
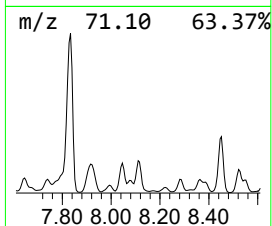
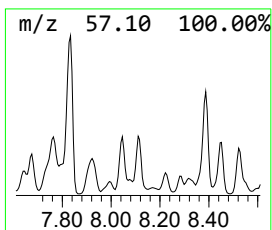
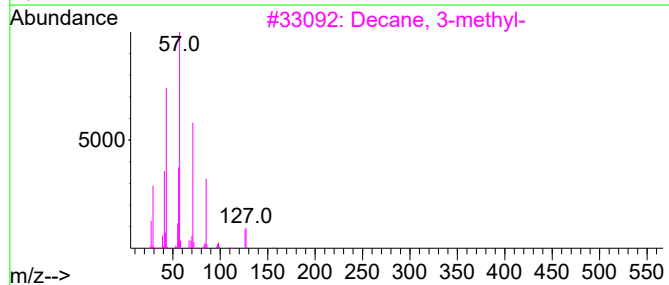
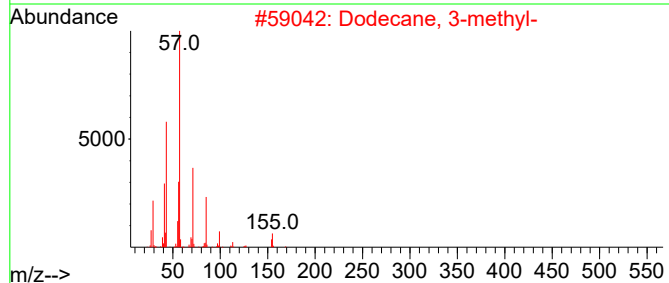
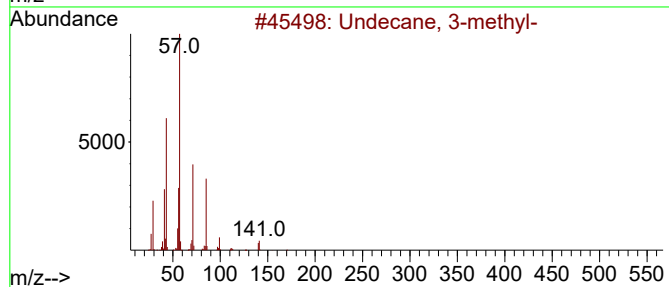
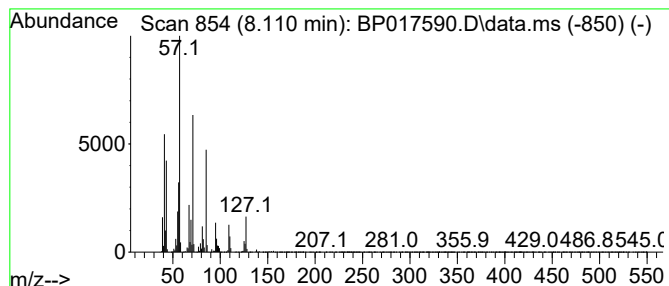
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 34 (DEL) Alkane: Straight-Chai... Concentration Rank 52

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.110	11.90 ng/ul	3242890	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3-methyl-	170	C12H26	001002-43-3	50	
2	Dodecane, 3-methyl-	184	C13H28	017312-57-1	50	
3	Decane, 3-methyl-	156	C11H24	013151-34-3	49	
4	Octane, 3,4,5,6-tetramethyl-	170	C12H26	062185-21-1	47	
5	10-Methylnonadecane	282	C20H42	056862-62-5	47	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

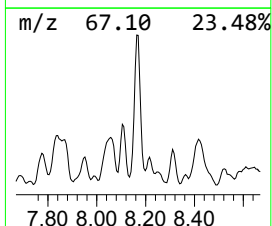
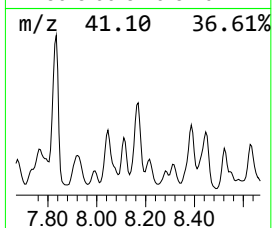
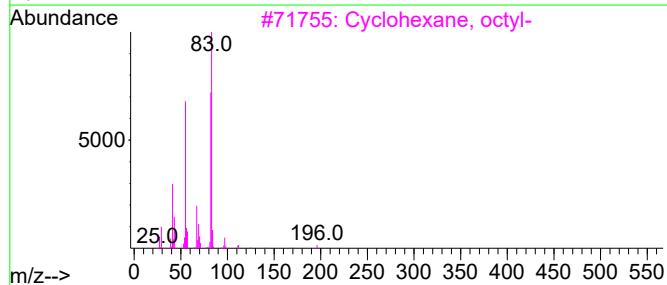
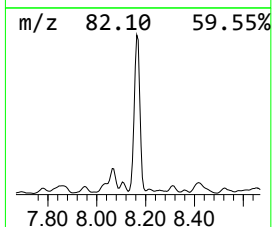
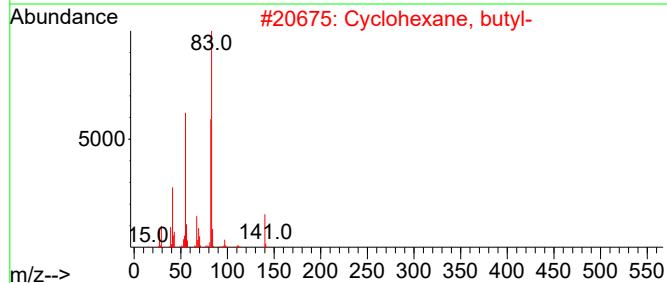
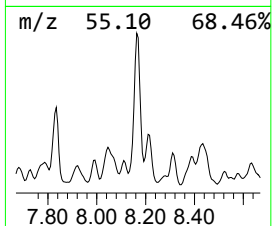
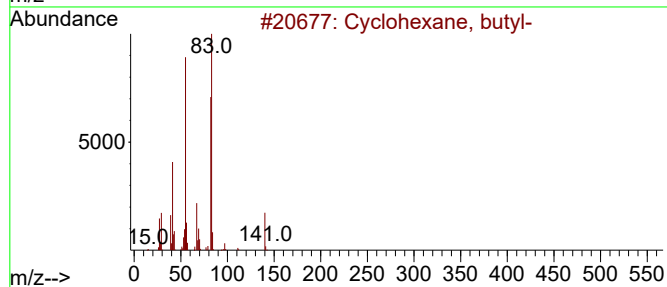
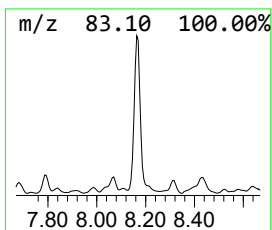
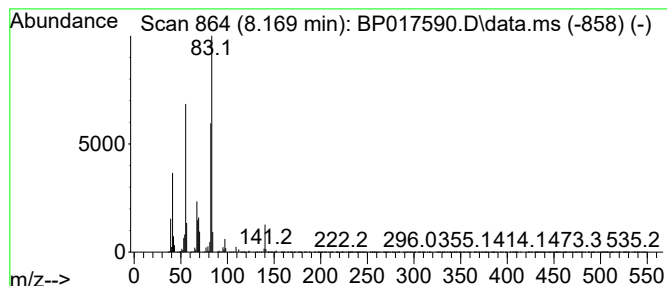
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 35 (DEL) Alkane: Cyclic8.169 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.169	35.07 ng/ul	9552070	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	90
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	87
3			Cyclohexane, octyl-	196	C14H28	001795-15-9	86
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	76
5			Cyclohexane, octyl-	196	C14H28	001795-15-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

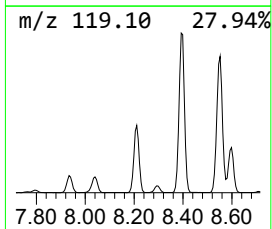
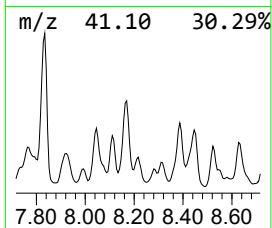
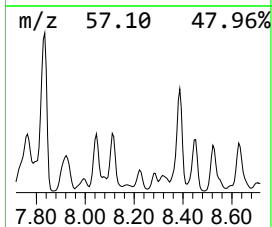
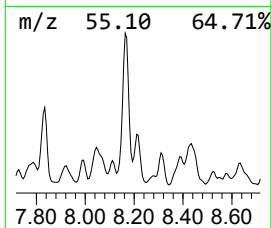
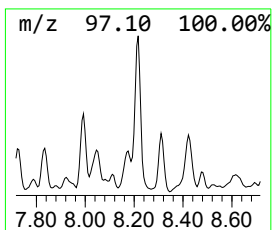
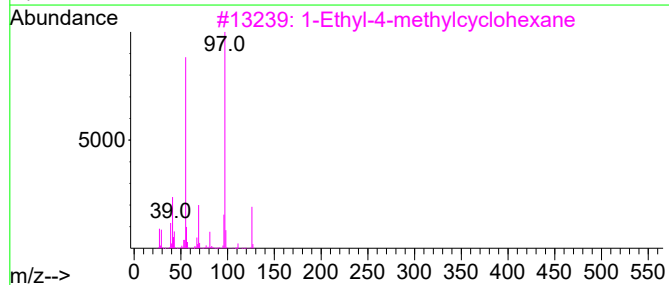
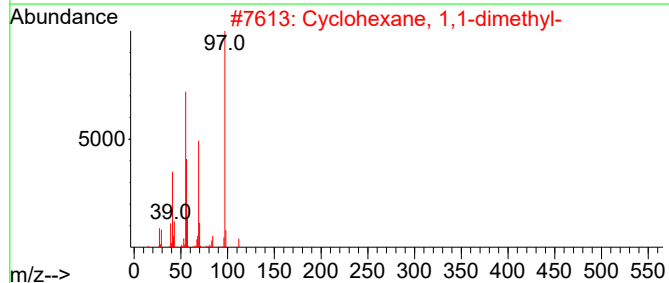
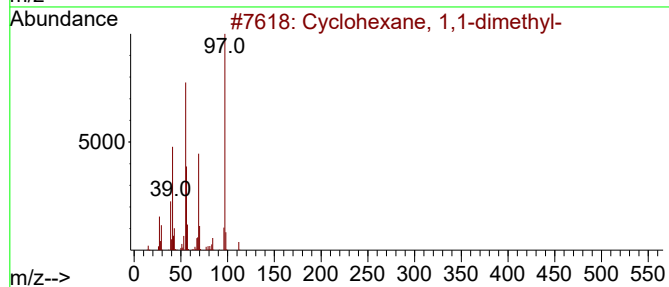
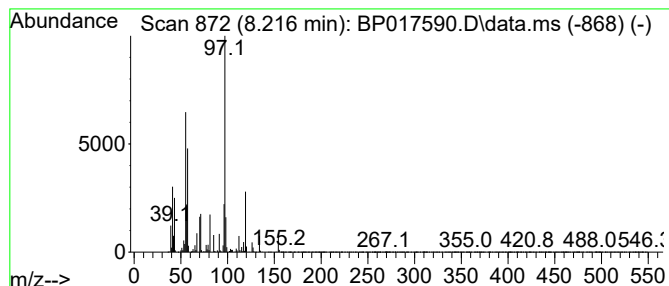
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 36 (DEL) Alkane: Cyclic8.216 Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.216	13.40 ng/ul	3650610	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	53
2	Cyclohexane, 1,1-dimethyl-	112	C8H16	000590-66-9	50
3	1-Ethyl-4-methylcyclohexane	126	C9H18	003728-56-1	50
4	Sulfurous acid, cyclohexylmethyl...	430	C25H50O3S	1000309-22-6	47
5	Cyclopentane, 1,1'-ethylidenebis-	166	C12H22	004413-21-2	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

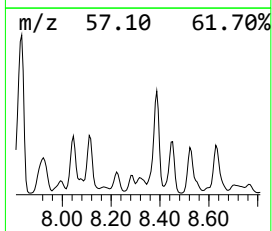
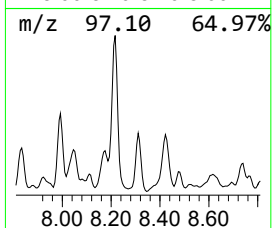
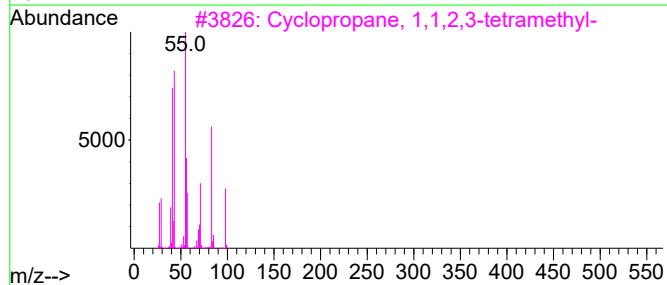
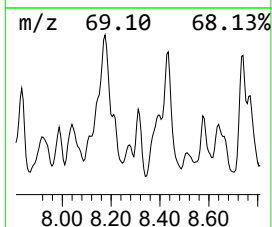
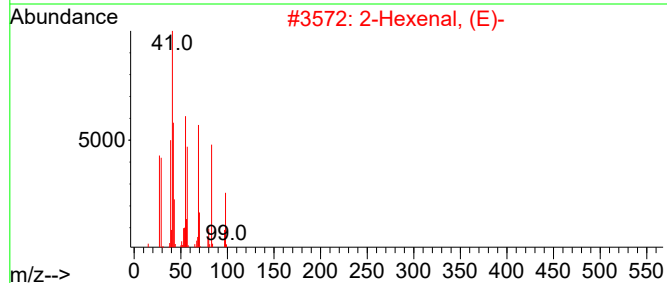
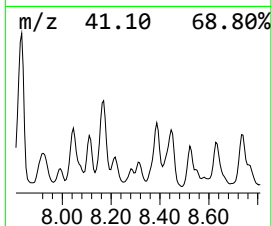
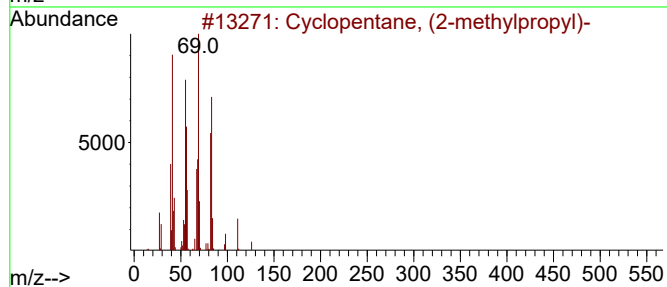
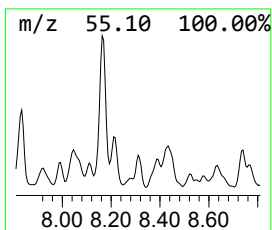
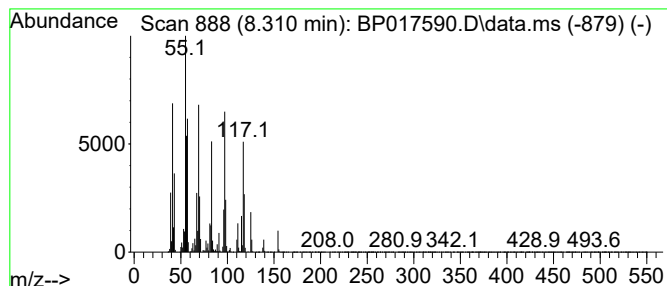
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 37 (DEL) Alkane: Cyclic8.310 Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.310	18.82 ng/ul	5125280	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclopentane, (2-methylpropyl)-	126	C9H18	003788-32-7	38
2			2-Hexenal, (E)-	98	C6H10O	006728-26-3	35
3			Cyclopropane, 1,1,2,3-tetramethyl-	98	C7H14	074752-93-5	30
4			Cycloheptane, methyl-	112	C8H16	004126-78-7	25
5			3-Heptene	98	C7H14	000592-78-9	15



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

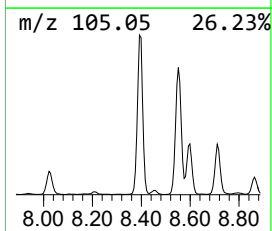
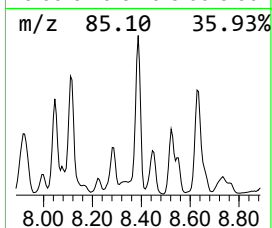
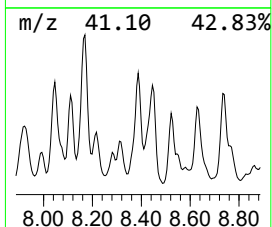
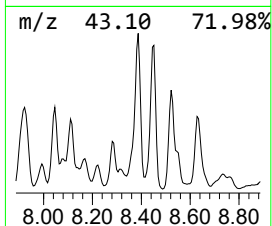
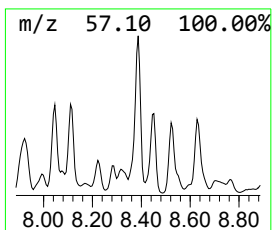
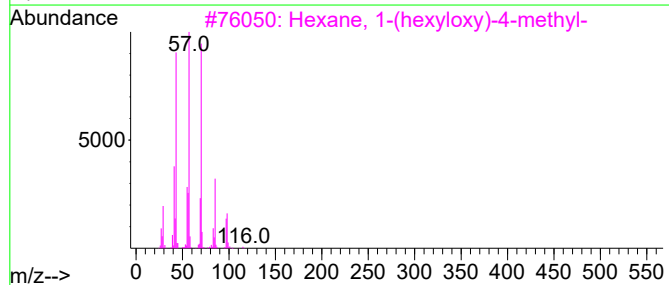
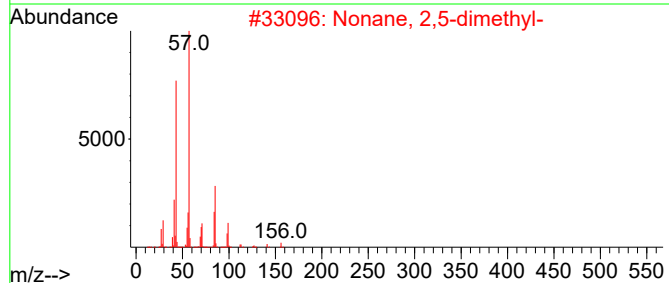
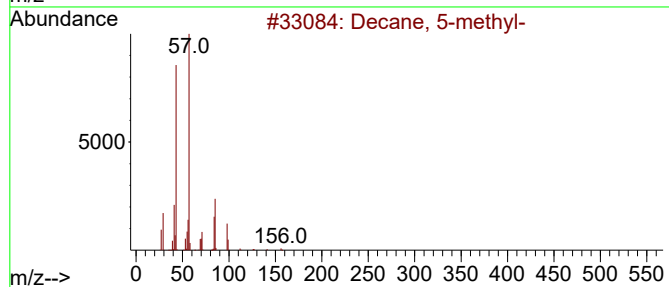
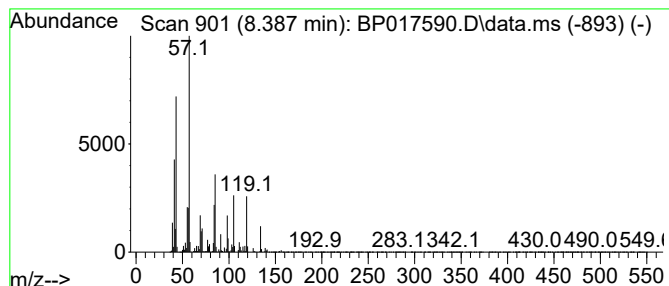
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.387	39.55 ng/ul	10773900	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Decane, 5-methyl-	156	C11H24	013151-35-4	49
2		Nonane, 2,5-dimethyl-	156	C11H24	017302-27-1	43
3		Hexane, 1-(hexyloxy)-4-methyl-	200	C13H28O	074421-20-8	38
4		Heptane, 4-ethyl-	128	C9H20	002216-32-2	38
5		Heptane, 4-ethyl-	128	C9H20	002216-32-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

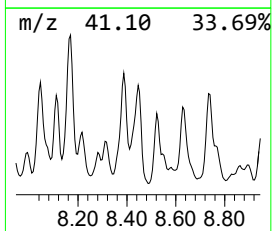
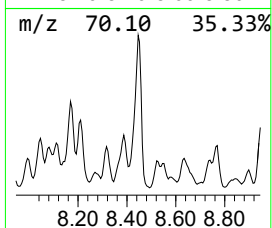
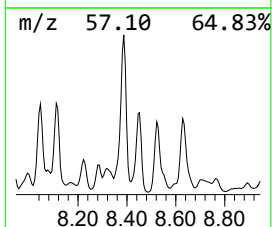
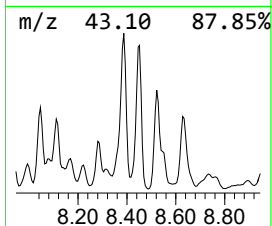
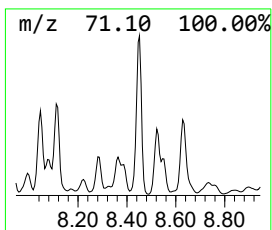
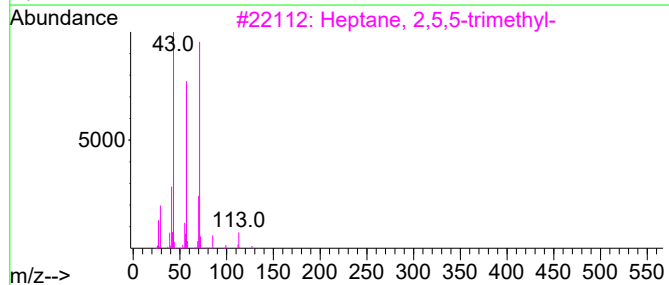
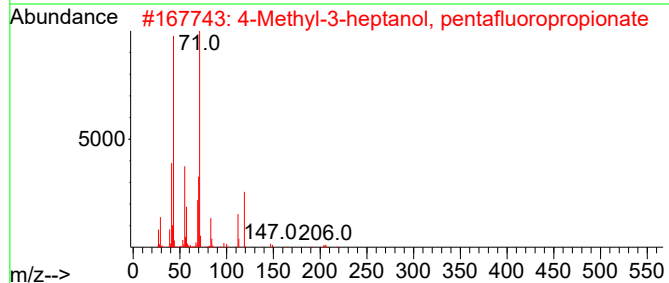
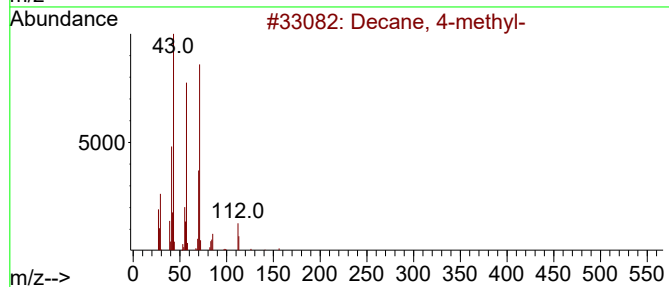
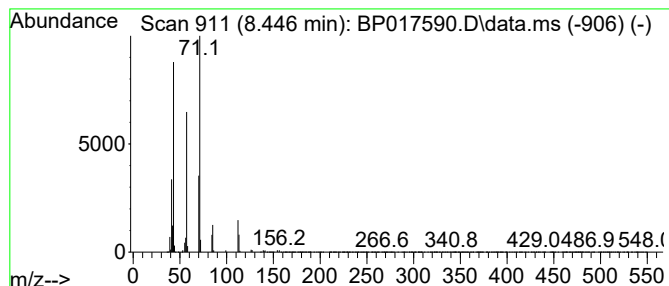
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 39 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.446	33.11 ng/ul	9018730	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	91
2			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	78
3			Heptane, 2,5,5-trimethyl-	142	C10H22	001189-99-7	78
4			Octane, 3-ethyl-	142	C10H22	005881-17-4	72
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

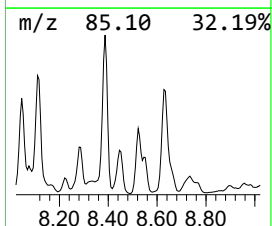
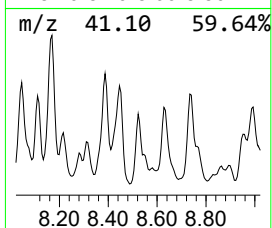
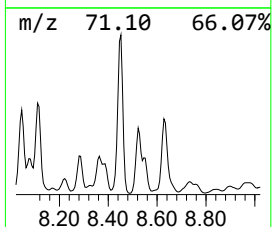
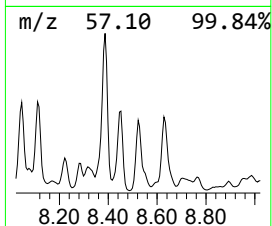
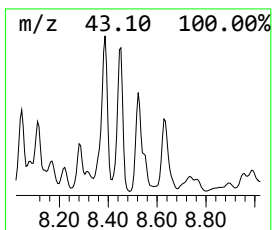
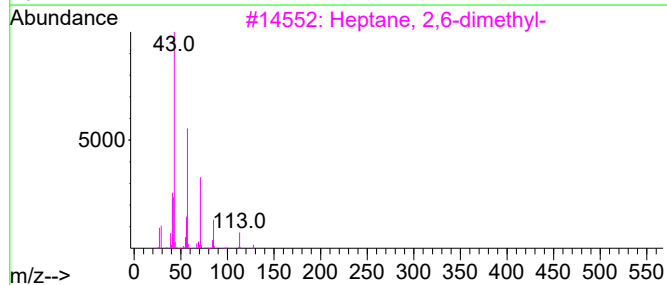
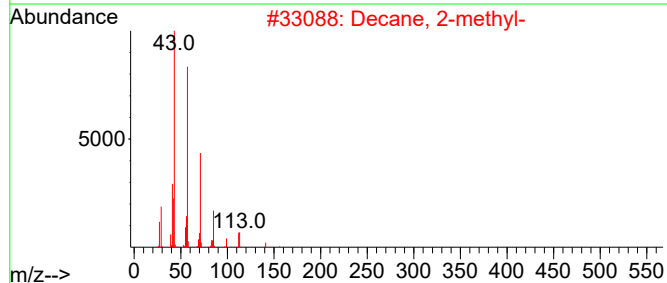
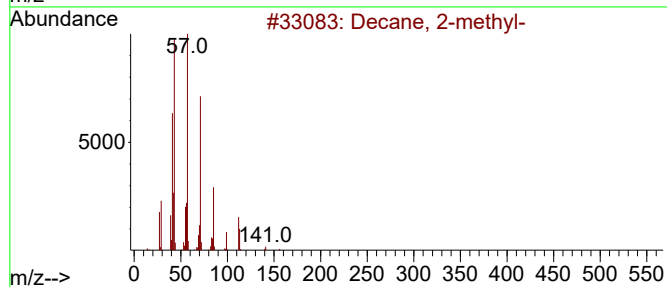
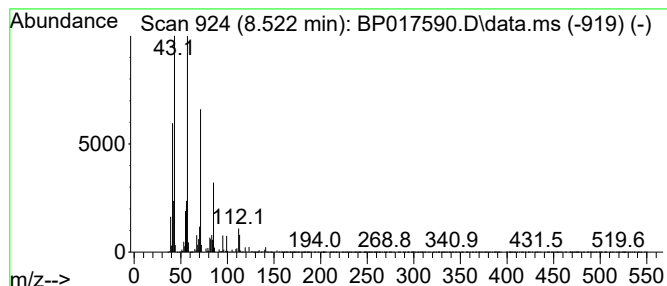
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.522	14.30 ng/ul	3894810	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 2-methyl-		156	C11H24	006975-98-0	86
2	Decane, 2-methyl-		156	C11H24	006975-98-0	81
3	Heptane, 2,6-dimethyl-		128	C9H20	001072-05-5	74
4	Hexadecane		226	C16H34	000544-76-3	59
5	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

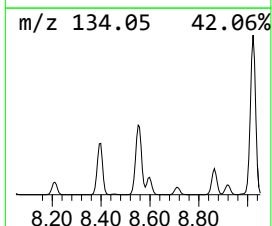
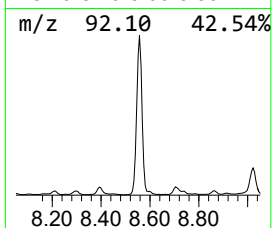
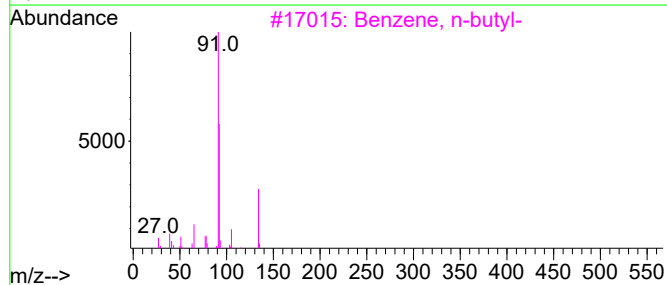
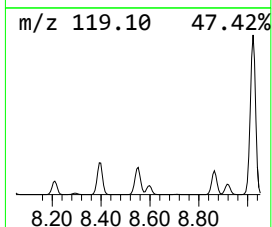
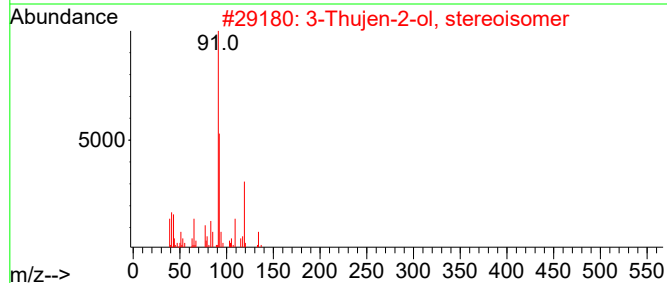
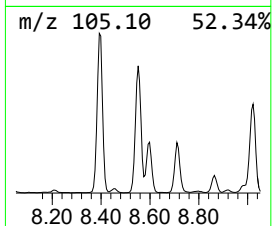
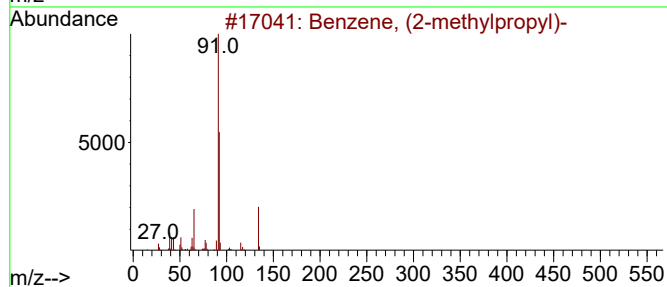
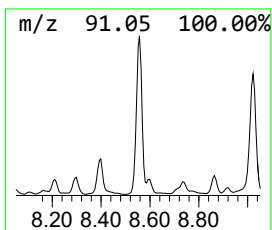
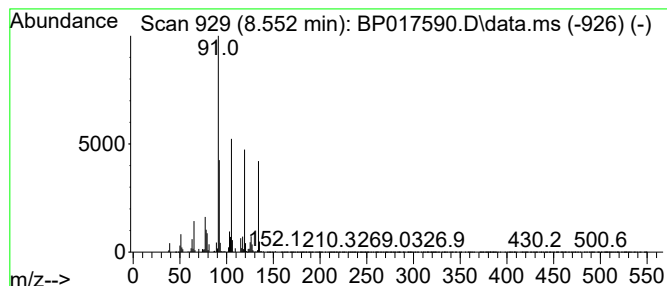
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 41 Benzene, (2-methylpropyl)- Concentration Rank 54

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.552	11.53 ng/ul	3139650	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (2-methylpropyl)-	134	C10H14	000538-93-2	53
2			3-Thujen-2-ol, stereoisomer	152	C10H16O	003310-03-0	50
3			Benzene, n-butyl-	134	C10H14	000104-51-8	49
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	38
5			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

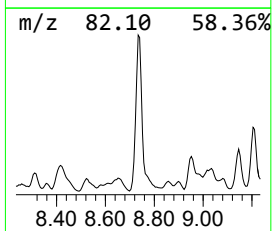
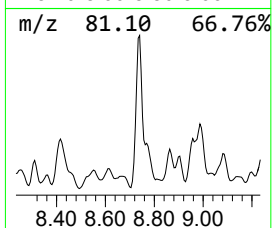
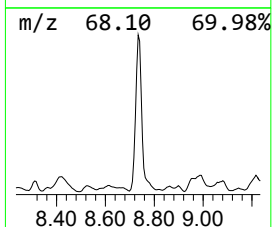
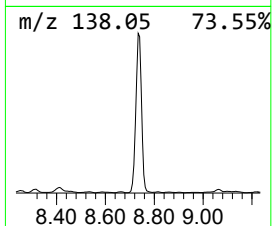
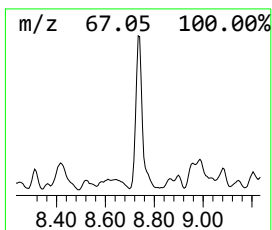
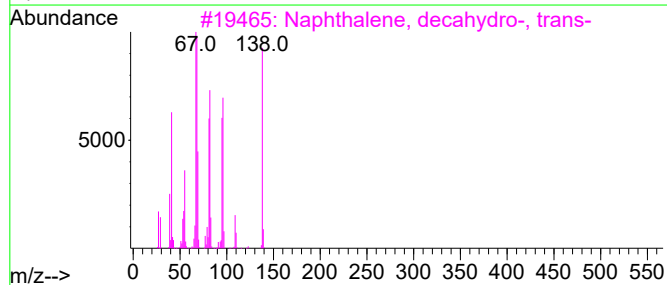
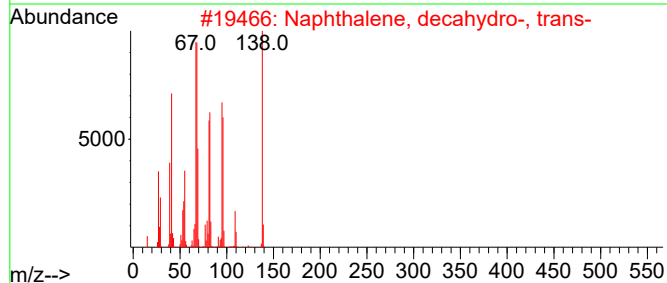
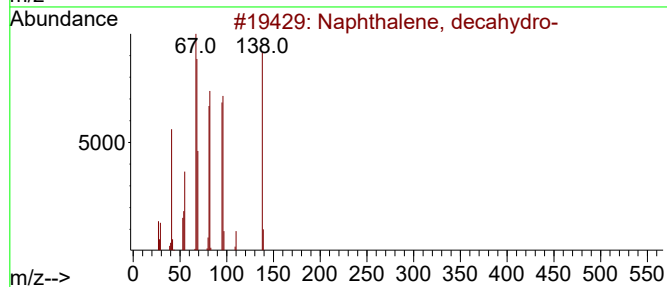
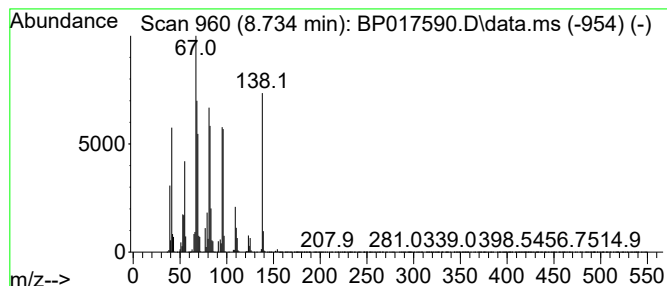
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 42 Naphthalene, decahydro- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.734	36.31 ng/ul	9891930	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, decahydro-	138	C10H18	000091-17-8	98
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	97
3			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
4			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	96
5			Naphthalene, decahydro-	138	C10H18	000091-17-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

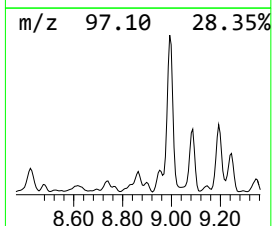
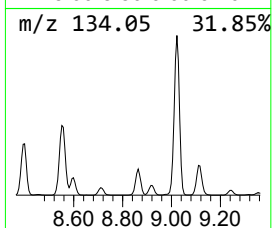
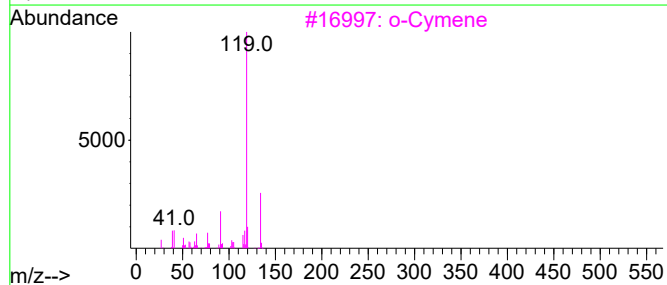
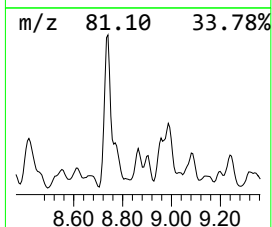
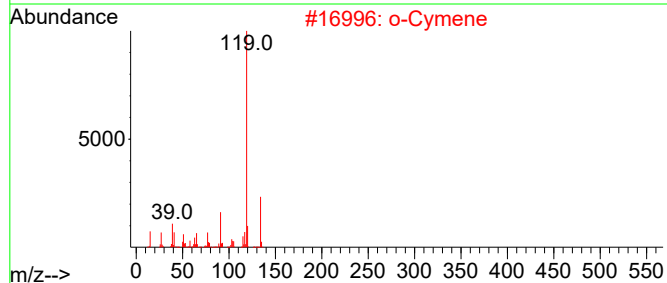
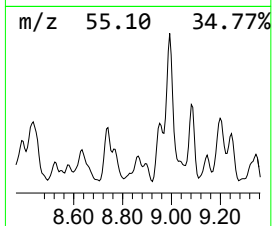
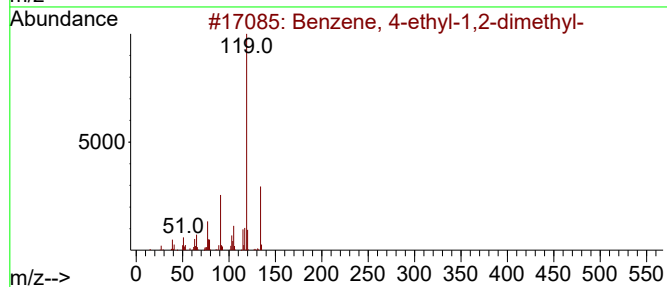
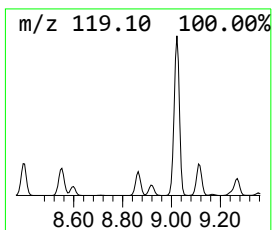
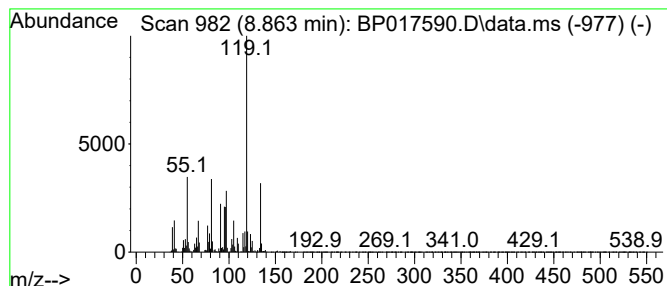
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 43 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 76

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.863	6.96 ng/ul	1896420	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			o-Cymene	134	C10H14	000527-84-4	93
3			o-Cymene	134	C10H14	000527-84-4	93
4			p-Cymene	134	C10H14	000099-87-6	93
5			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

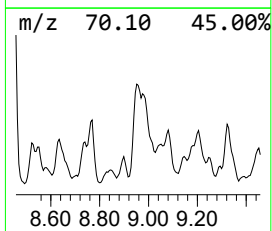
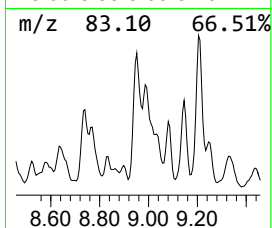
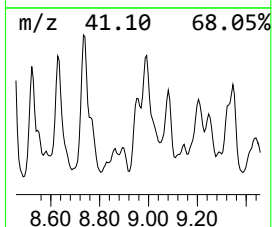
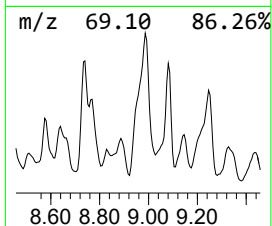
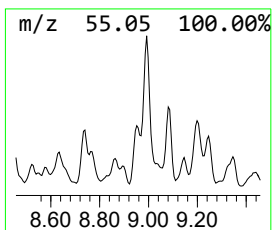
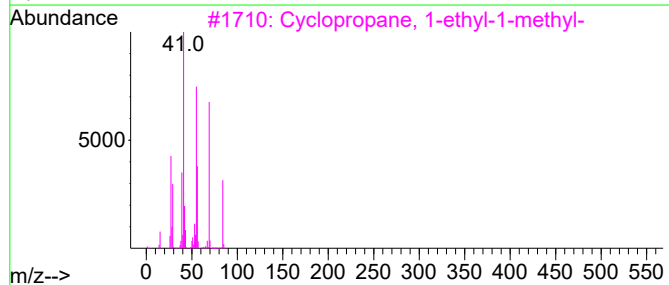
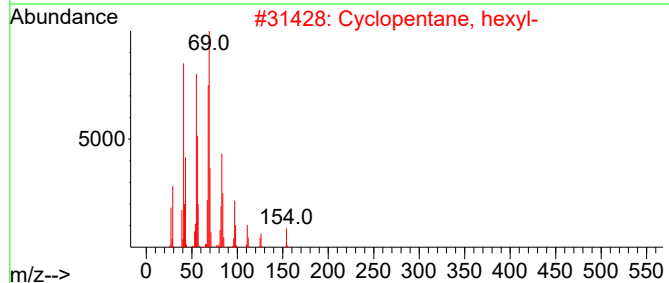
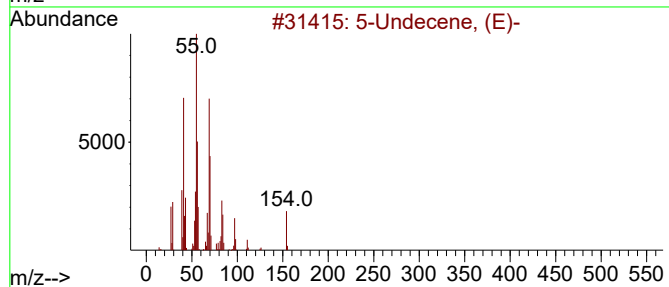
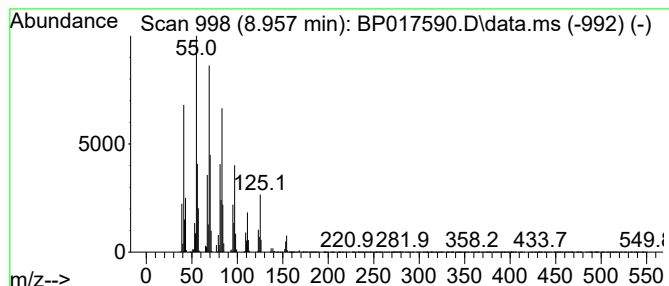
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.957	15.14 ng/ul	4123380	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Undecene, (E)-	154	C11H22	000764-97-6	58
2	Cyclopentane, hexyl-	154	C11H22	004457-00-5	49
3	Cyclopropane, 1-ethyl-1-methyl-	84	C6H12	053778-43-1	41
4	1-Hexene, 3,3-dimethyl-	112	C8H16	003404-77-1	38
5	3-Dodecene, (E)-	168	C12H24	007206-14-6	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

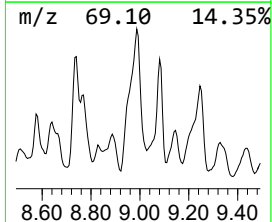
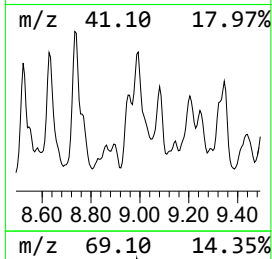
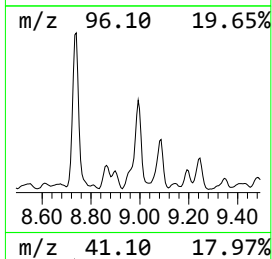
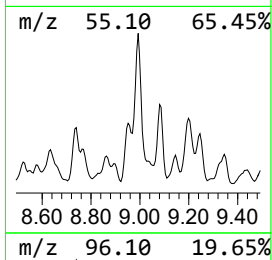
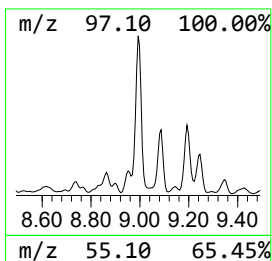
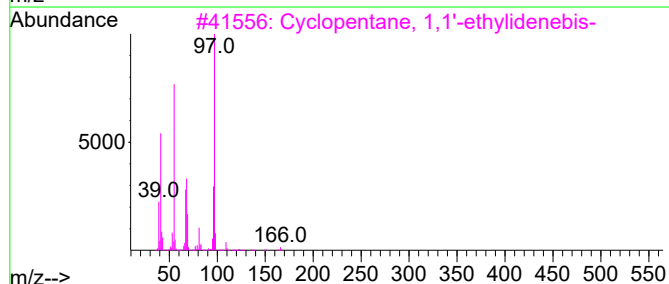
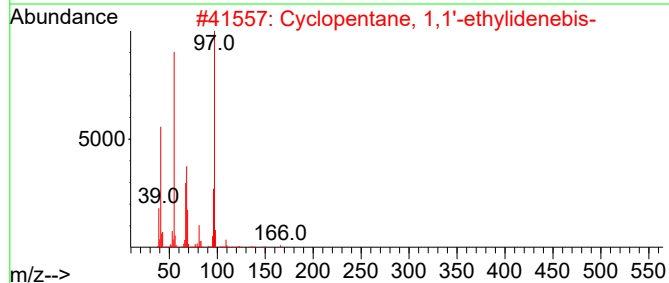
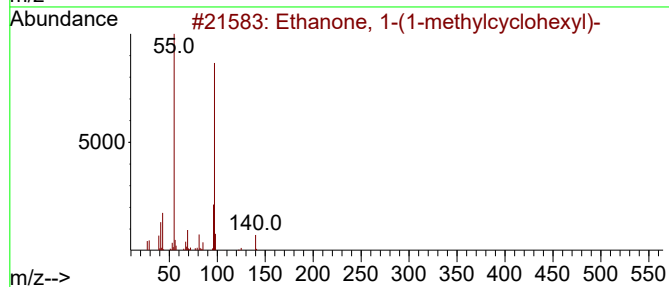
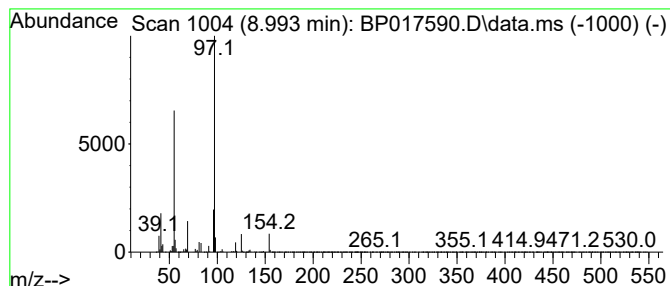
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 45 Ethanone, 1-(1-methylcyclohexyl) Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.993	23.97 ng/ul	6528970	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Ethanone, 1-(1-methylcyclohexyl)-		140	C9H16O	002890-62-2	72
2	Cyclopentane, 1,1'-ethylidenebis-		166	C12H22	004413-21-2	64
3	Cyclopentane, 1,1'-ethylidenebis-		166	C12H22	004413-21-2	64
4	Hept-2-ene, 2,4,4,6-tetramethyl-		154	C11H22	103982-58-7	53
5	Cycloheptane, bromo-		176	C7H13Br	002404-35-5	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

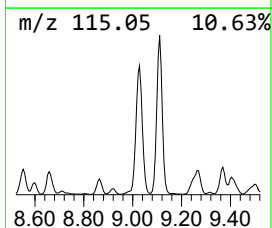
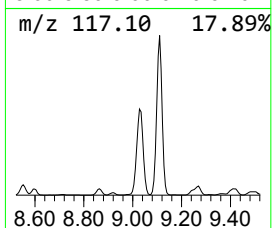
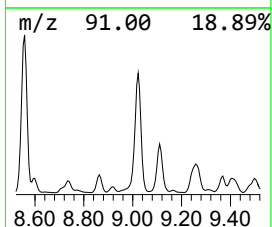
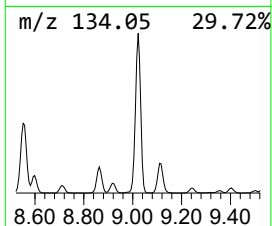
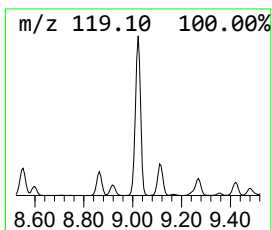
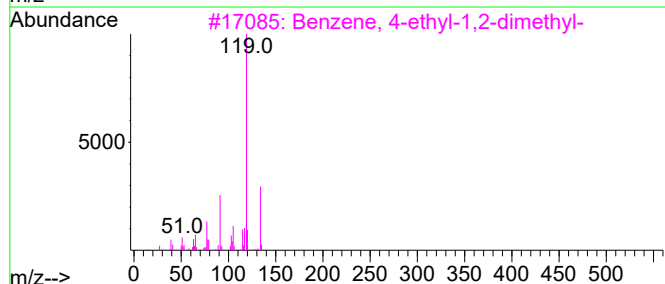
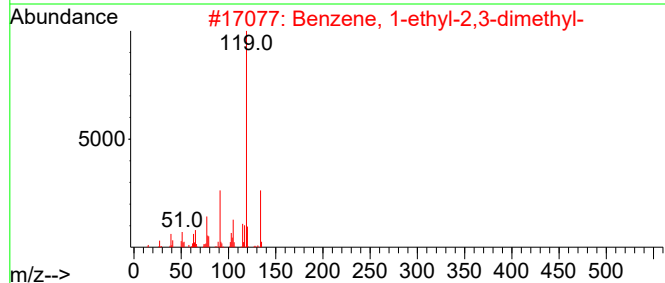
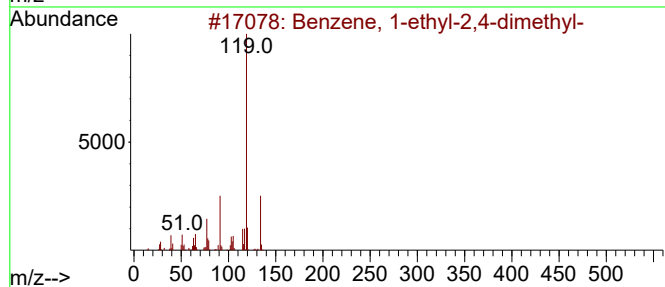
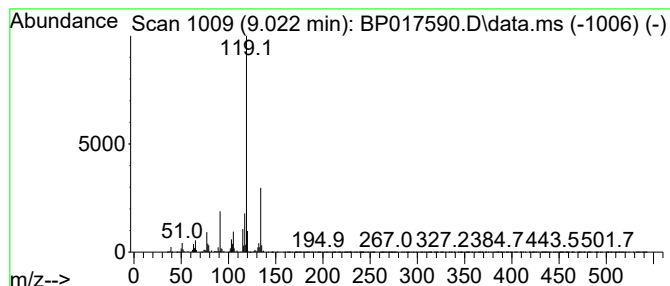
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 46 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.022	30.64 ng/ul	8346760	1,4-Dichlorobenzene-d4	7.940

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
2		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	95
3		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4		o-Cymene	134	C10H14	000527-84-4	94
5		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

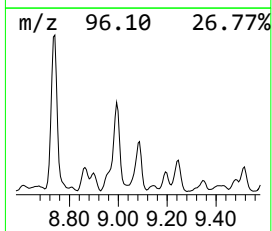
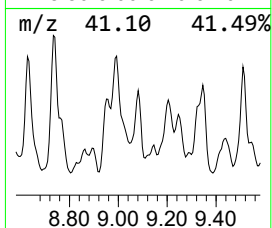
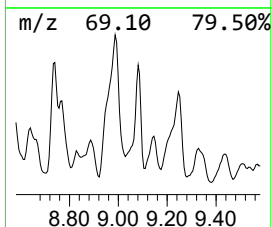
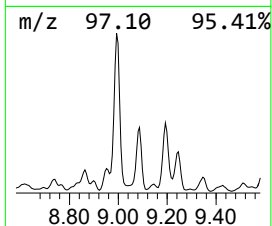
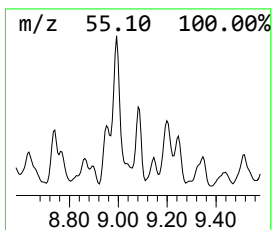
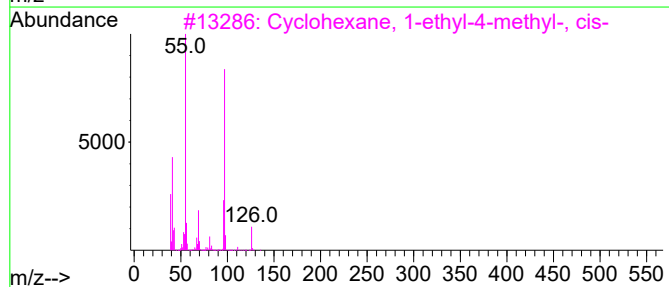
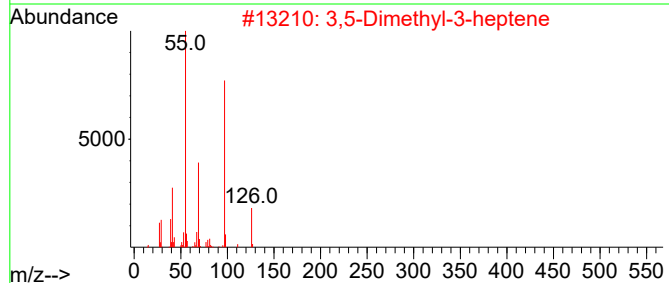
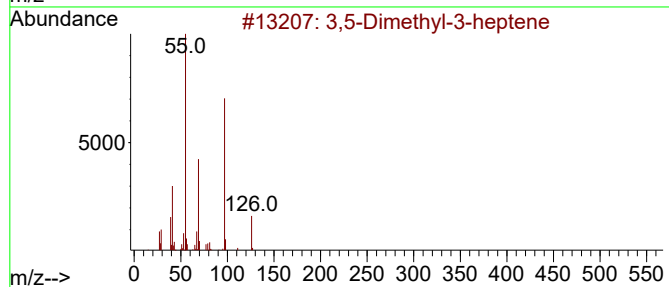
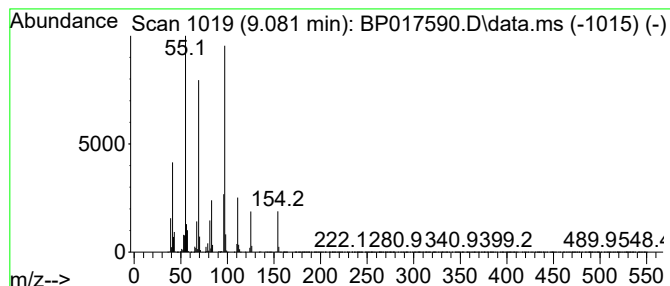
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.081	12.26 ng/ul	3339150	1,4-Dichlorobenzene-d4	7.940	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	53
2	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	53
3	Cyclohexane, 1-ethyl-4-methyl-, ...	126	C9H18	004926-78-7	50
4	2(5H)-Furanone, 5,5-dimethyl-	112	C6H8O2	020019-64-1	47
5	Cyclopentanecarboxylic acid, eth...	140	C8H12O2	016523-06-1	47



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

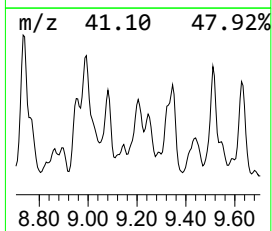
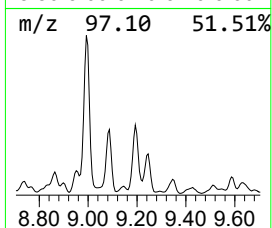
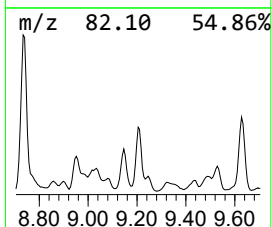
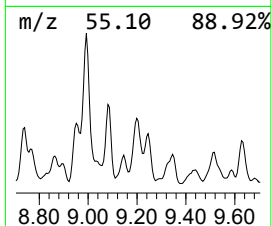
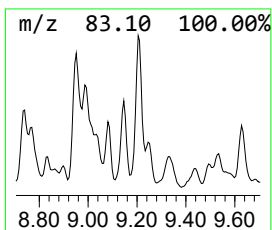
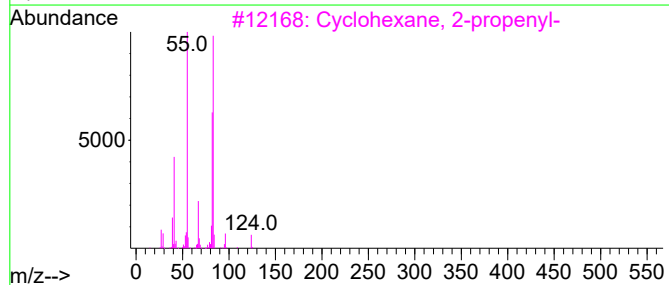
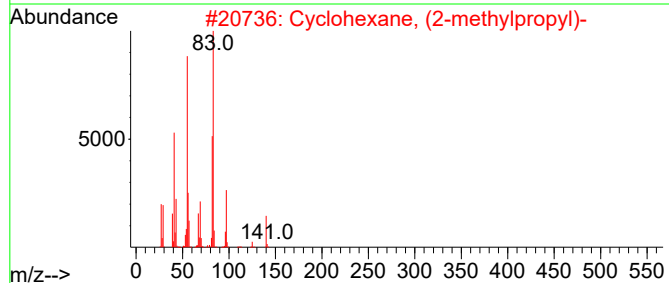
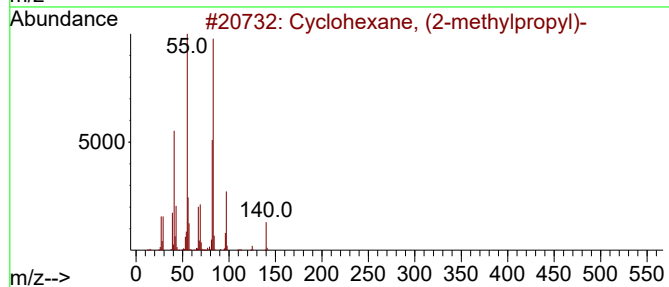
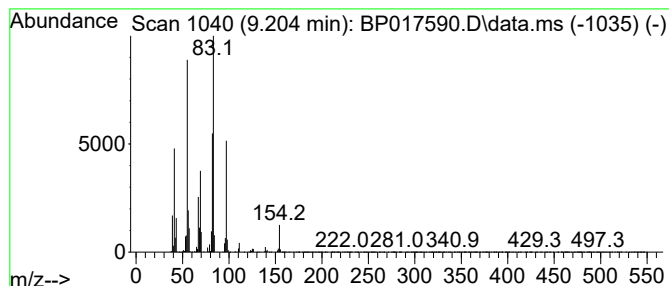
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 48 (DEL) Alkane: Cyclic9.204 Concentration Rank 71

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.204	8.23 ng/ul	2240530	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	64
2			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	64
3			Cyclohexane, 2-propenyl-	124	C9H16	002114-42-3	58
4			Cyclohexane, octyl-	196	C14H28	001795-15-9	53
5			Cyclohexane, butyl-	140	C10H20	001678-93-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

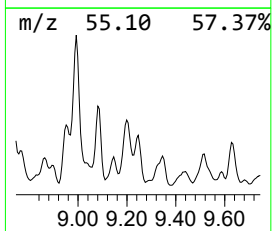
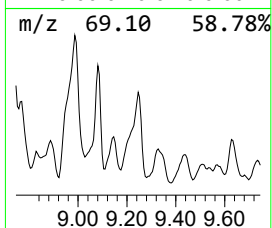
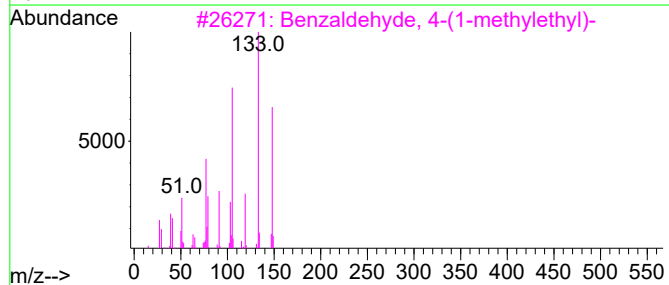
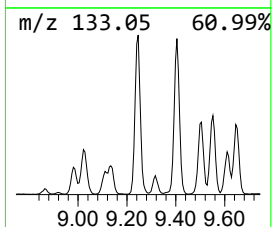
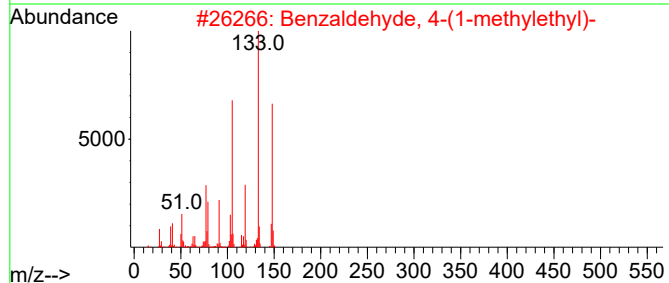
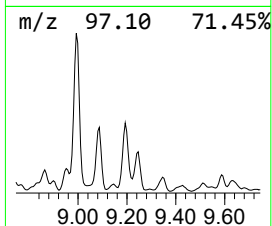
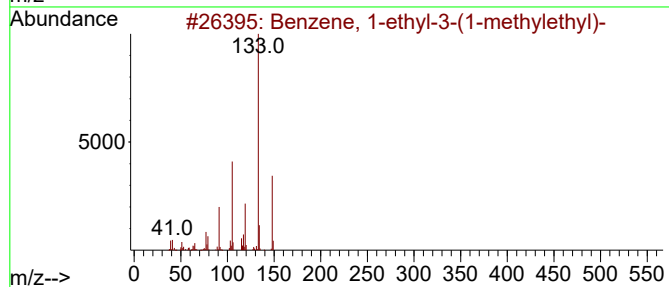
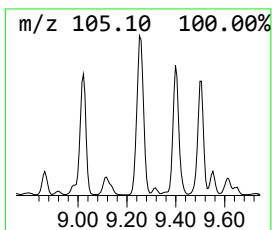
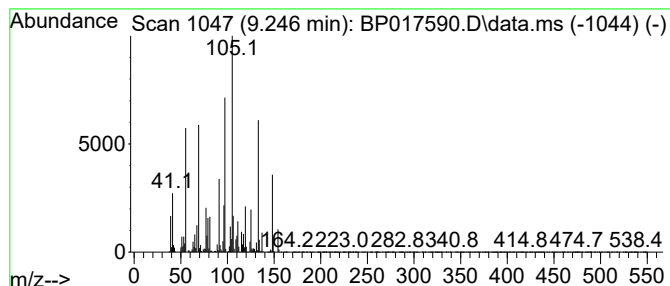
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 49 Benzene, 1-ethyl-3-(1-methy... Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.246	13.75 ng/ul	3746630	1,4-Dichlorobenzene-d4	7.940

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	64
2			Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	46
3			Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	41
4			Benzaldehyde, 4-(1-methylethyl)-	148	C10H12O	000122-03-2	41
5			3,4-Dimethylcumene	148	C11H16	1000370-34-1	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

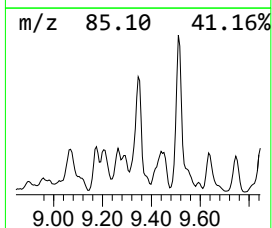
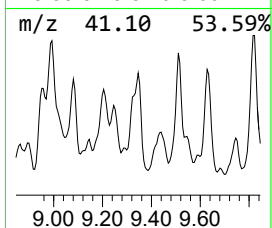
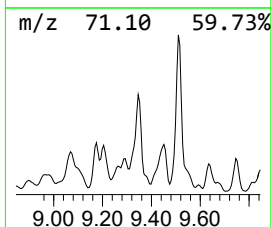
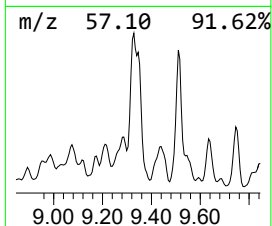
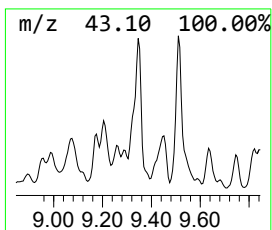
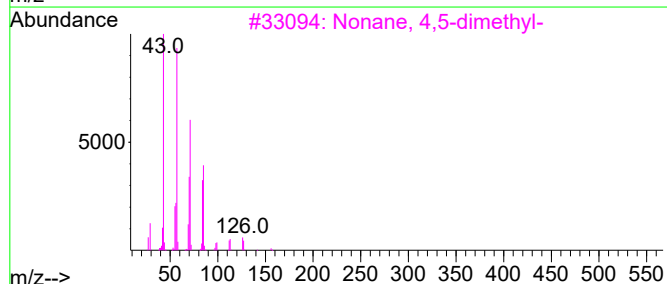
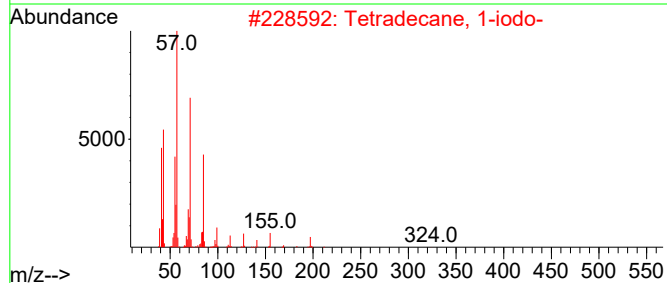
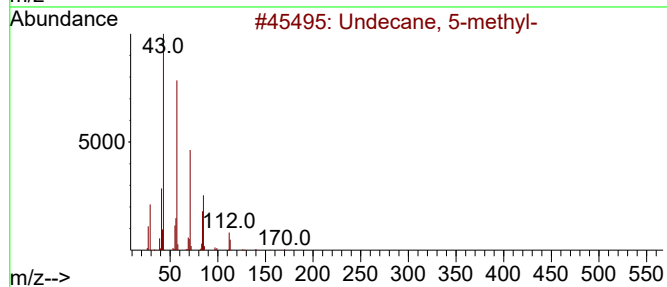
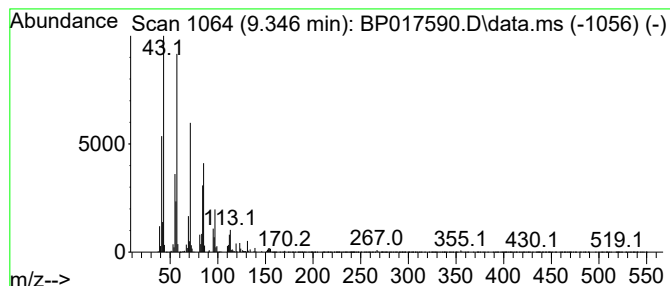
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD			R.T.
9.346	20.09 ng/ul	5471930	1,4-Dichlorobenzene-d4			7.940
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane, 5-methyl-		170	C12H26	001632-70-8	81
2	Tetradecane, 1-iodo-		324	C14H29I	019218-94-1	70
3	Nonane, 4,5-dimethyl-		156	C11H24	017302-23-7	59
4	Decane, 3,8-dimethyl-		170	C12H26	017312-55-9	58
5	Octane, 2,3,3-trimethyl-		156	C11H24	062016-30-2	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

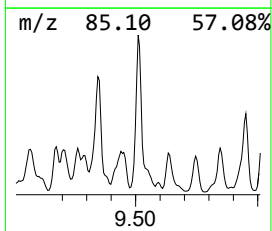
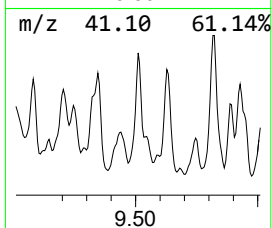
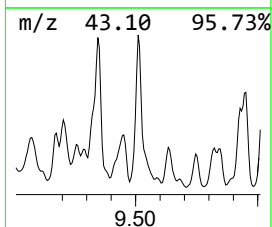
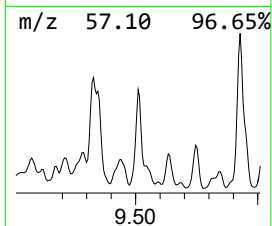
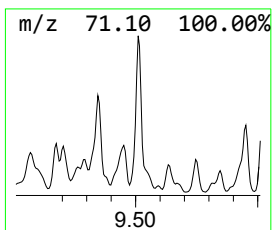
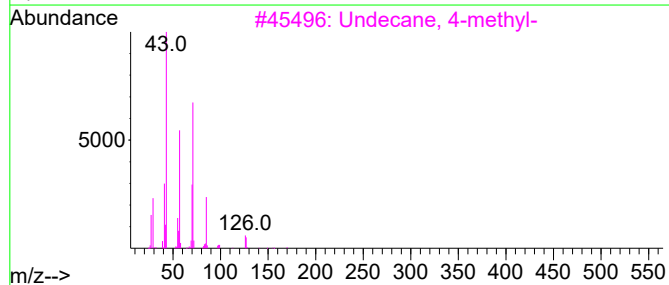
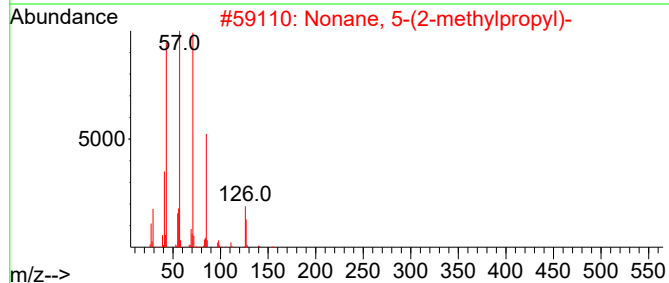
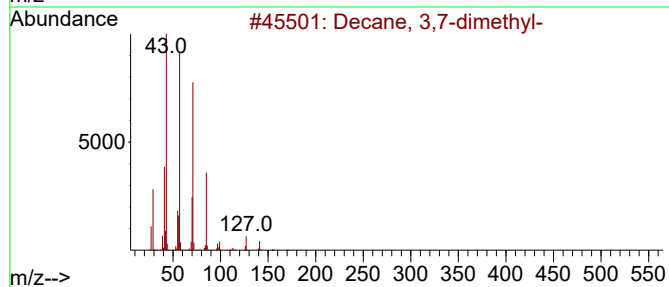
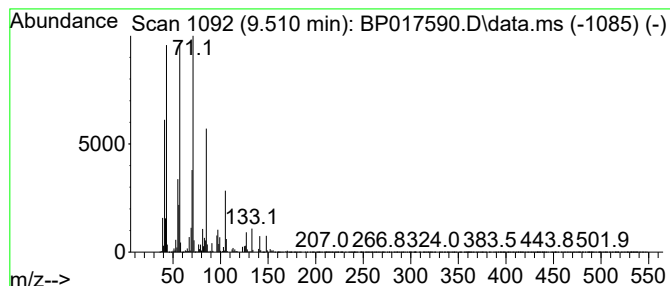
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 51 (DEL) Alkane: Straight-Chai... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.510	26.59 ng/ul	5231370	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	70
2			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	53
3			Undecane, 4-methyl-	170	C12H26	002980-69-0	52
4			Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	46
5			Undecane, 2,3-dimethyl-	184	C13H28	017312-77-5	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

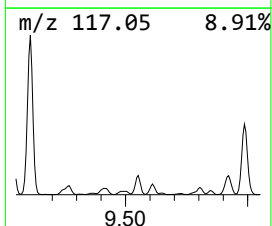
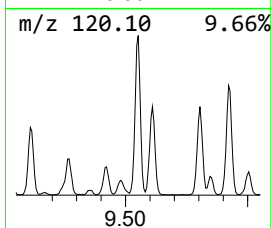
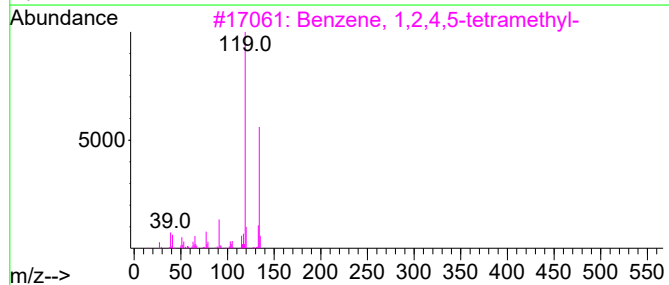
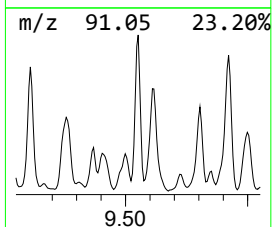
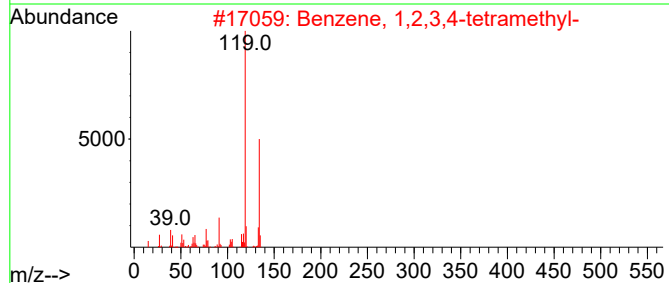
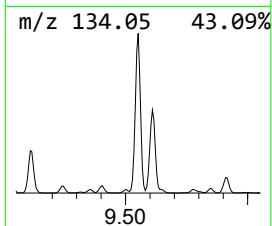
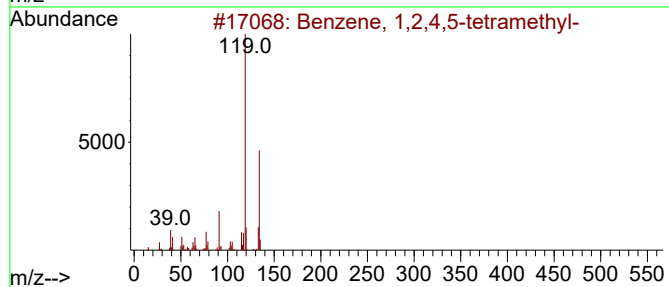
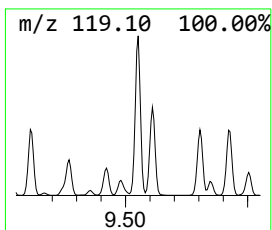
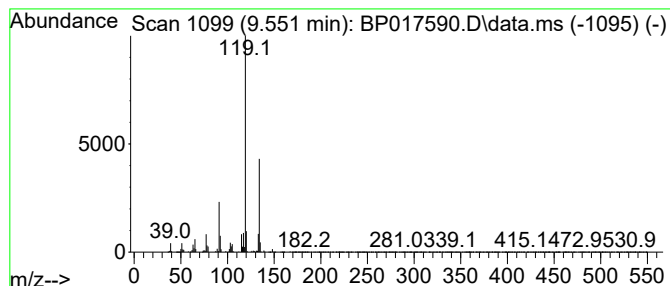
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 52 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.551	19.13 ng/ul	3762490	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	94
3			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

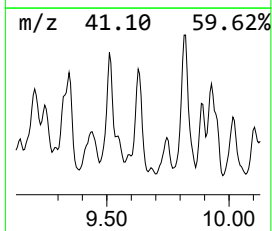
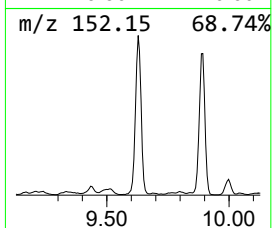
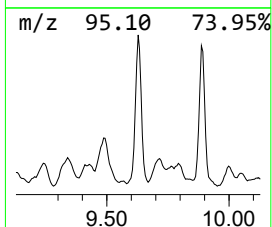
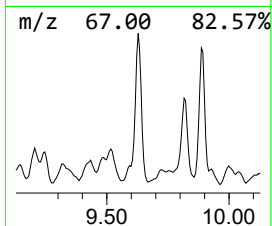
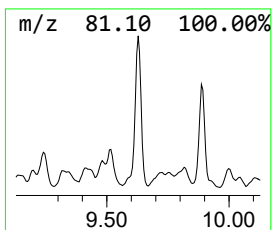
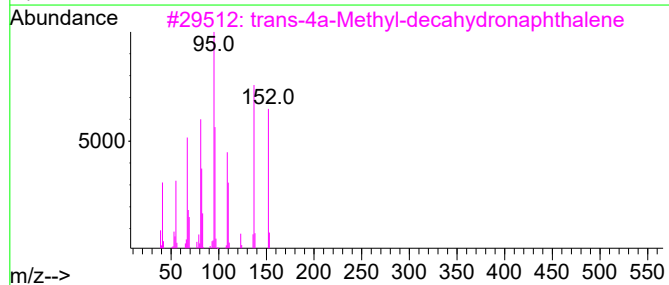
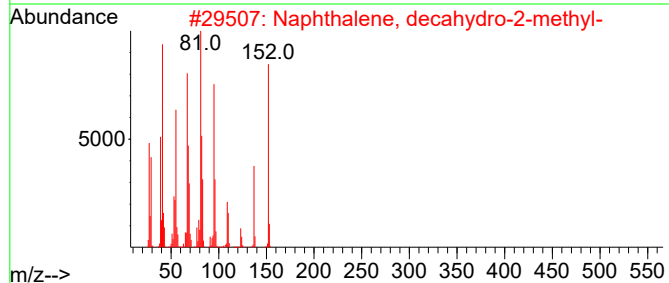
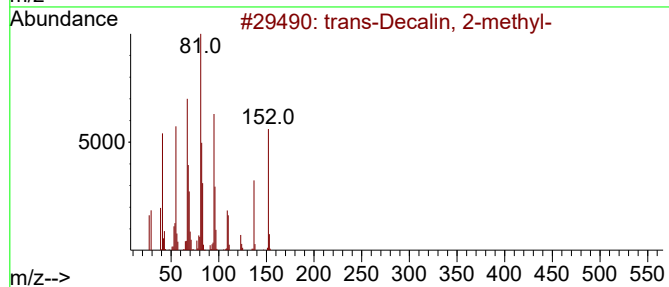
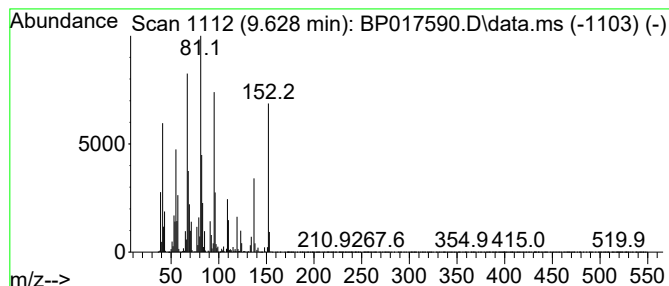
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 53 trans-Decalin, 2-methyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.628	46.71 ng/ul	9188160	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-Decalin, 2-methyl-	152	C11H20	1000152-47-3	95
2			Naphthalene, decahydro-2-methyl-	152	C11H20	002958-76-1	94
3			trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	89
4			trans-4a-Methyl-decahydronaphtha...	152	C11H20	002547-27-5	86
5			Cyclohexanone, 5-methyl-2-(1-met...	152	C10H16O	015932-80-6	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

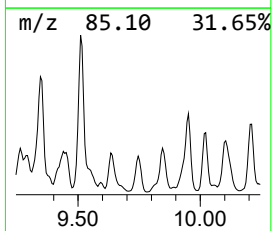
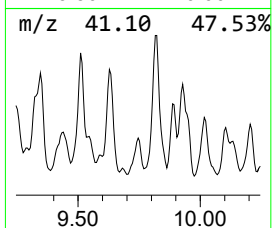
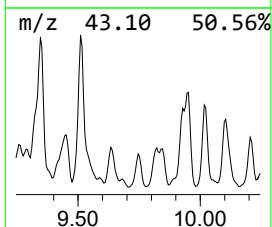
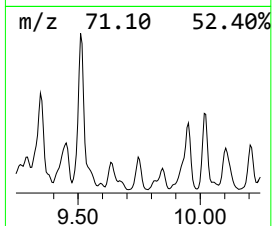
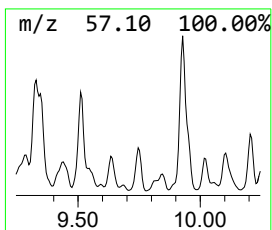
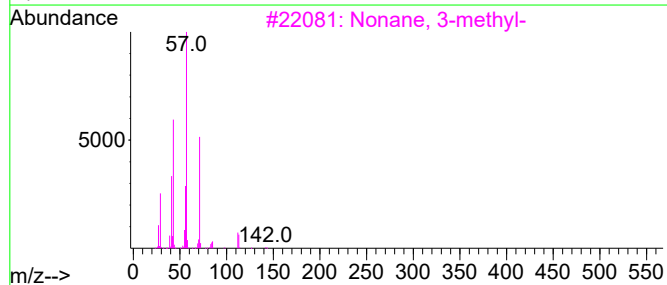
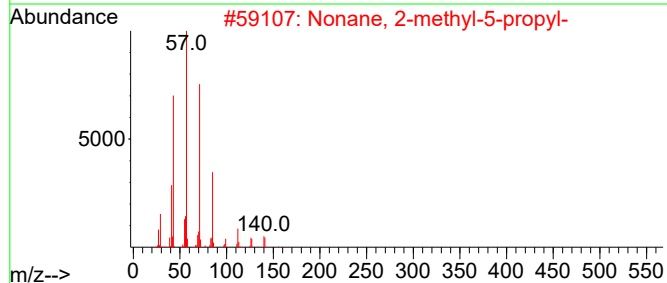
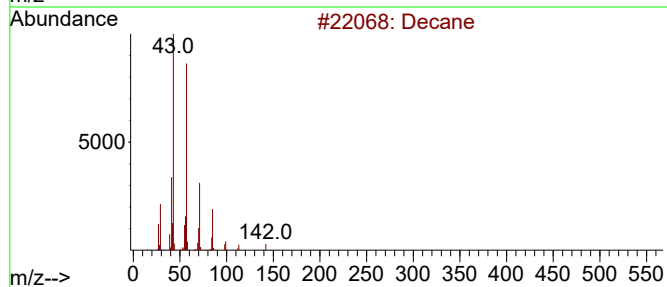
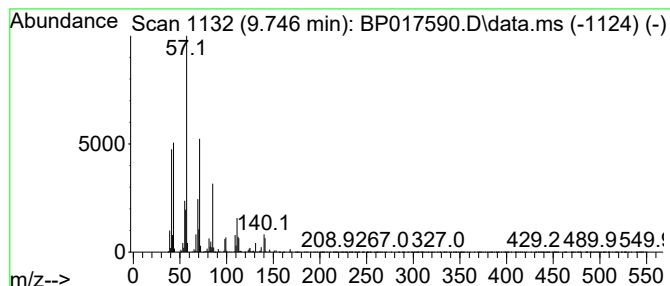
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.746	10.45 ng/ul	2055880	Naphthalene-d8	10.775	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane	142	C10H22	000124-18-5	76
2	Nonane, 2-methyl-5-propyl-	184	C13H28	031081-17-1	64
3	Nonane, 3-methyl-	142	C10H22	005911-04-6	64
4	Undecane, 3,7-dimethyl-	184	C13H28	017301-29-0	59
5	Undecane, 5-methyl-	170	C12H26	001632-70-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

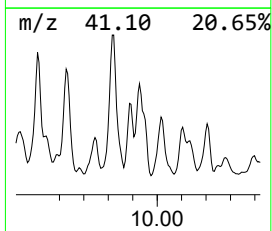
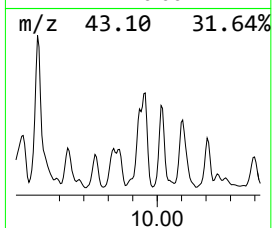
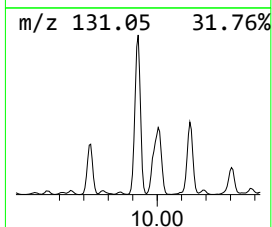
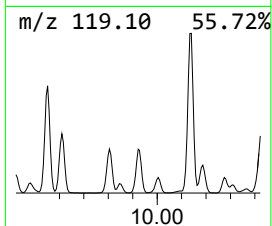
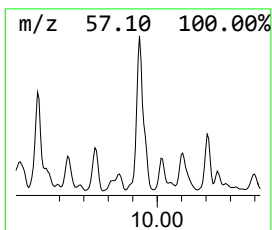
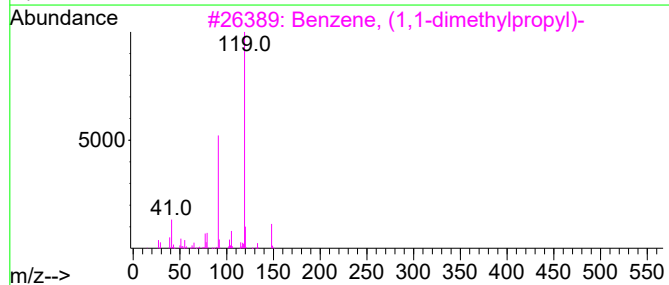
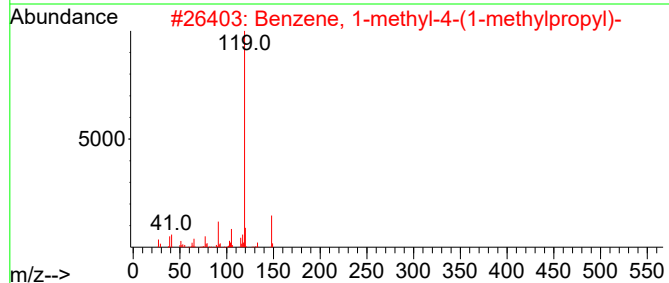
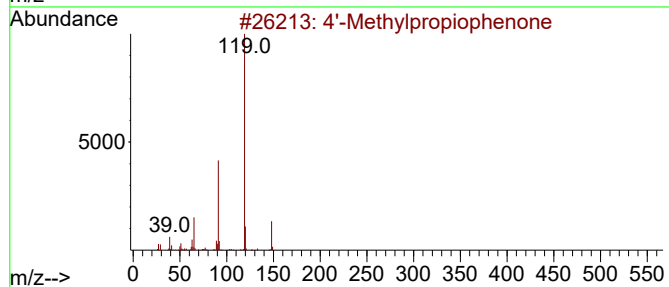
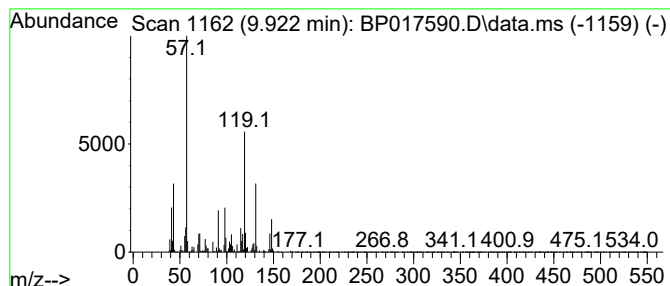
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 55 unknown-03 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
9.922	36.10 ng/ul	7101430	Naphthalene-d8	10.775	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4'-Methylpropiophenone	148	C10H12O	005337-93-9	42
2	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	30
3	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	30
4	4'-Methylpropiophenone	148	C10H12O	005337-93-9	27
5	Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

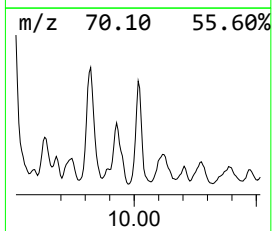
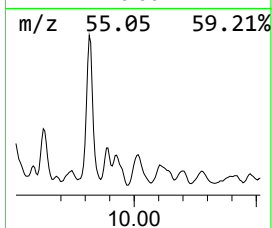
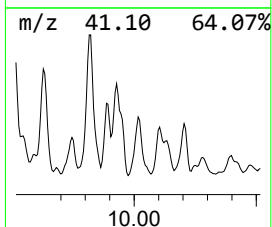
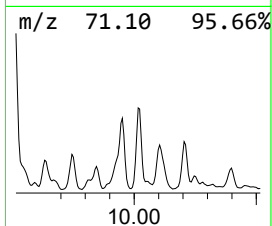
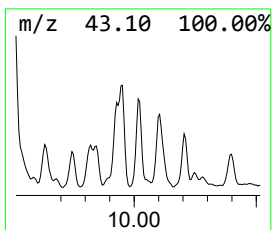
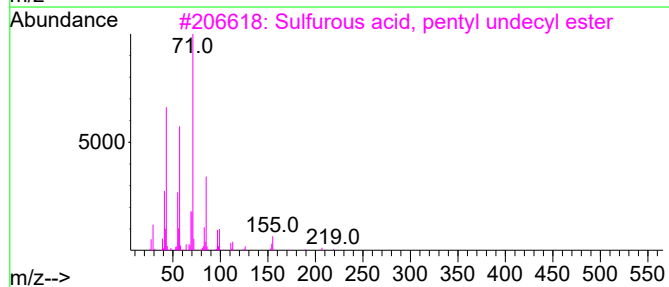
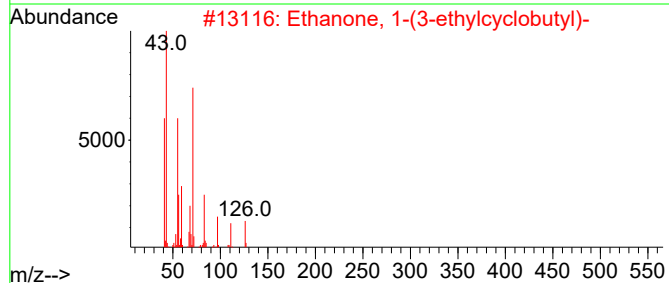
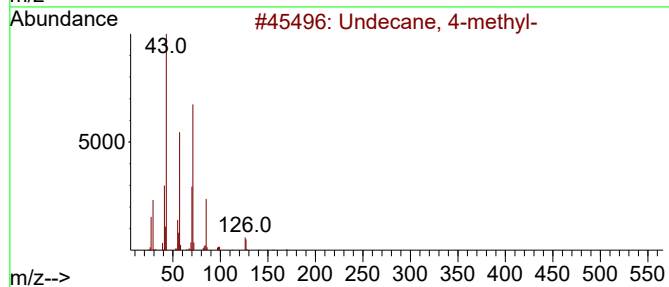
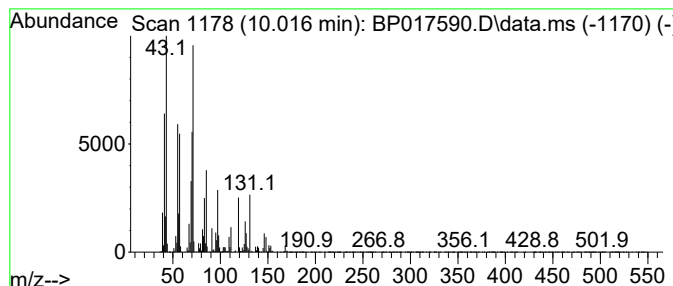
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 56 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.016	22.39 ng/ul	4404490	Naphthalene-d8	10.775	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 4-methyl-	170	C12H26	002980-69-0	53
2	Ethanone, 1-(3-ethylcyclobutyl)-	126	C8H14O	056335-71-8	43
3	Sulfurous acid, pentyl undecyl e...	306	C16H34O3S	1000309-14-4	38
4	4,4-Dimethyl-1-hexene	112	C8H16	001647-08-1	35
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

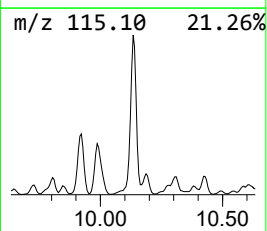
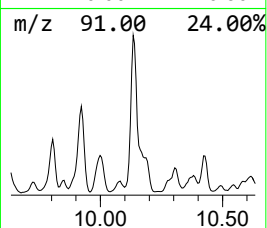
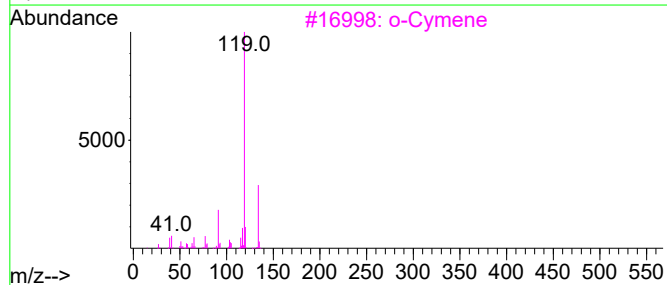
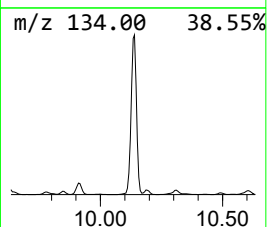
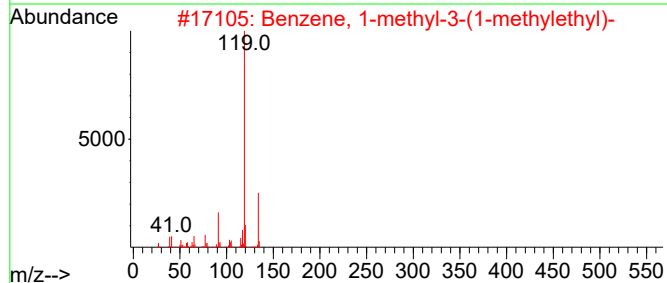
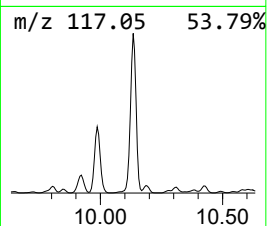
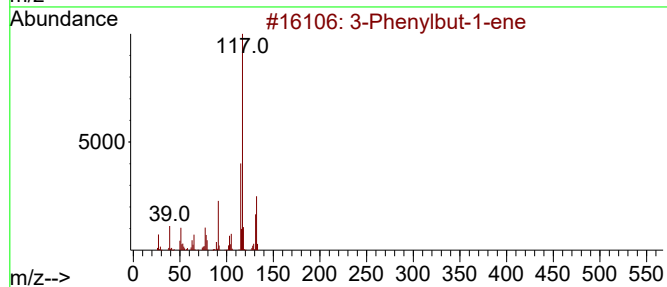
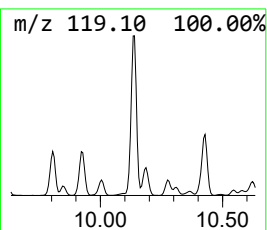
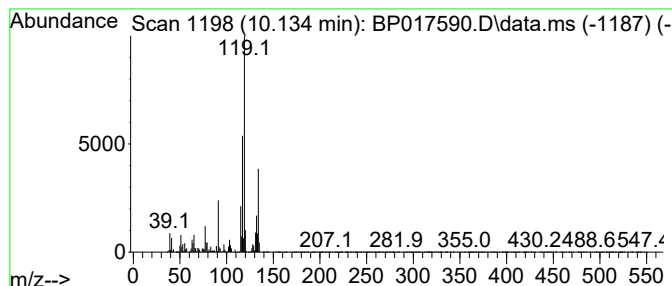
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 57 3-Phenylbut-1-ene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.134	50.86 ng/ul	10005300	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
2			Benzene, 1-methyl-3-(1-methylethyl)-	134	C10H14	000535-77-3	64
3			o-Cymene	134	C10H14	000527-84-4	64
4			p-Cymene	134	C10H14	000099-87-6	64
5			o-Cymene	134	C10H14	000527-84-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

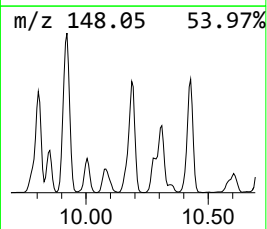
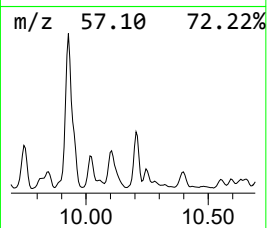
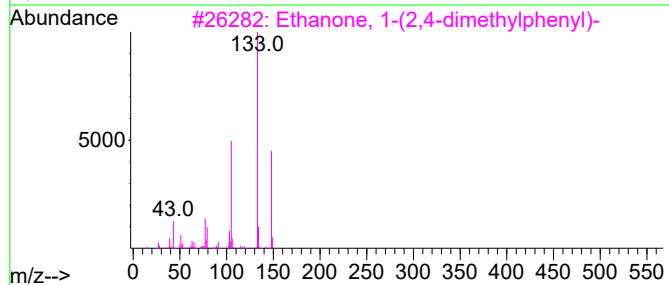
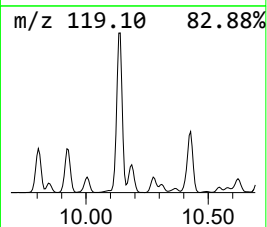
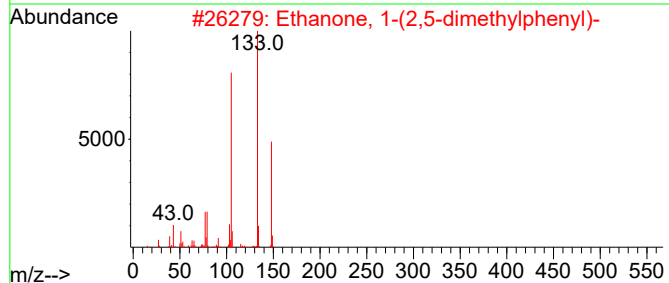
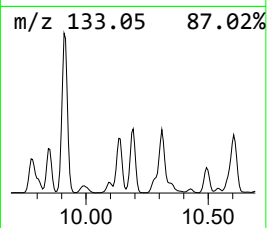
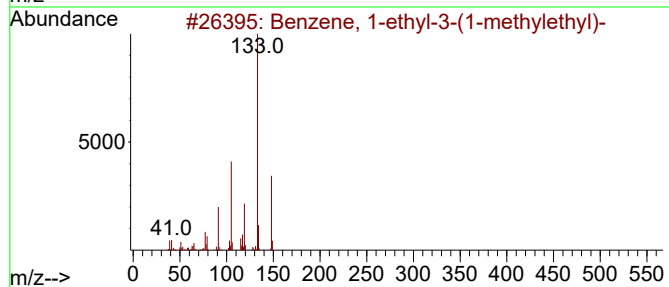
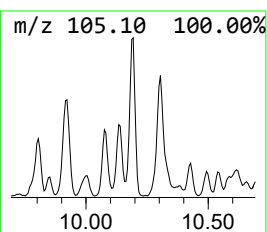
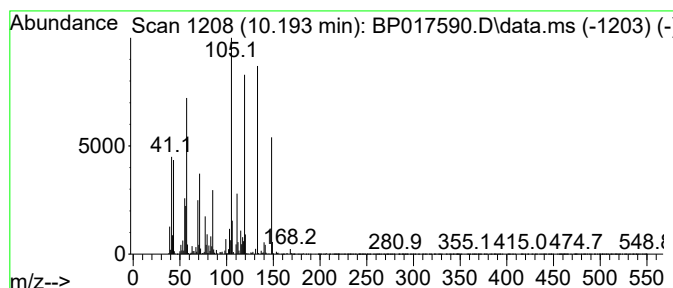
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 58 Ethanone, 1-(2,5-dimethylph... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.193	15.64 ng/ul	3076460	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3-(1-methylethyl)-	148	C11H16	004920-99-4	55
2			Ethanone, 1-(2,5-dimethylphenyl)-	148	C10H12O	002142-73-6	50
3			Ethanone, 1-(2,4-dimethylphenyl)-	148	C10H12O	000089-74-7	46
4			1-(2,3-Dimethylphenyl)ethanone	148	C10H12O	002142-71-4	38
5			Butyric acid, 2-phenyl-, dec-2-y...	304	C20H32O2	1000406-86-4	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

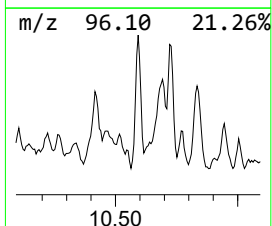
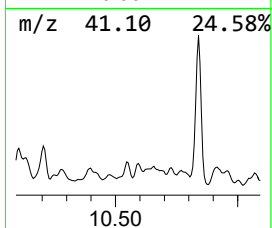
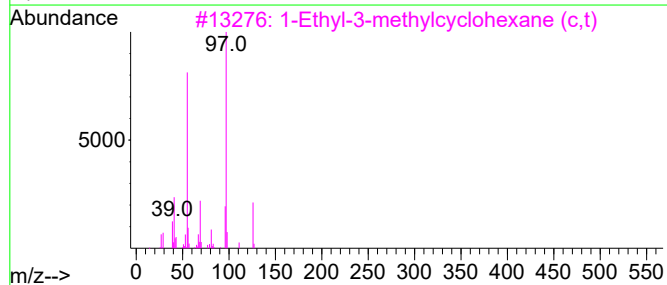
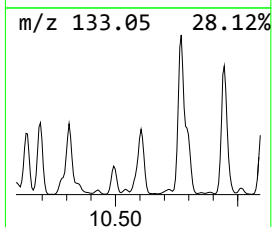
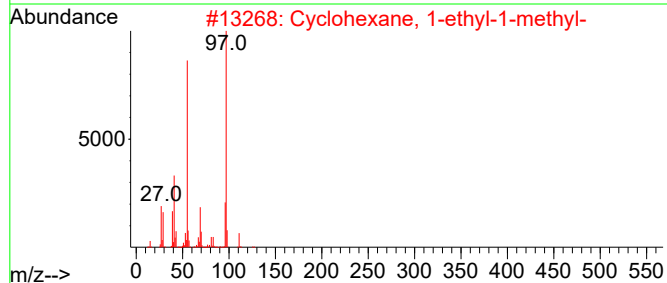
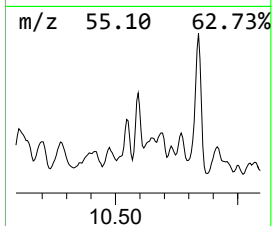
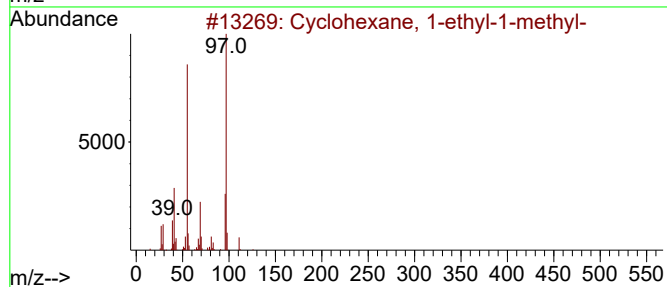
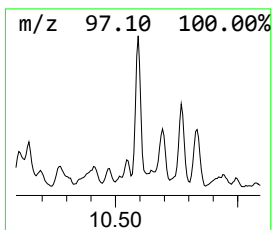
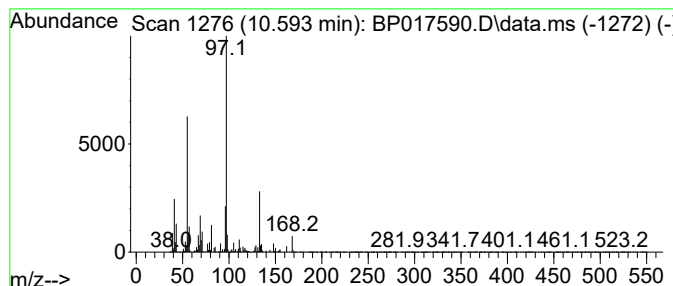
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 59 (DEL) Alkane: Cyclic10.593 Concentration Rank 67

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.593	8.58 ng/ul	1687450	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	70
2			Cyclohexane, 1-ethyl-1-methyl-	126	C9H18	004926-90-3	58
3			1-Ethyl-3-methylcyclohexane (c,t)	126	C9H18	003728-55-0	53
4			cis-1-Ethyl-3-methyl-cyclohexane	126	C9H18	019489-10-2	50
5			5-Amino-3-methylpyrazole	97	C4H7N3	031230-17-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

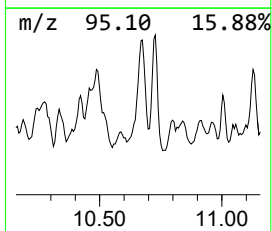
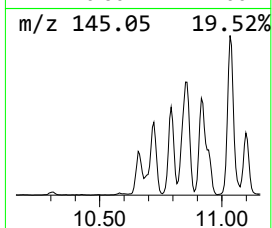
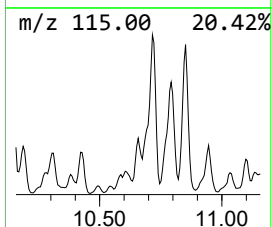
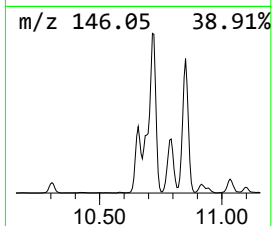
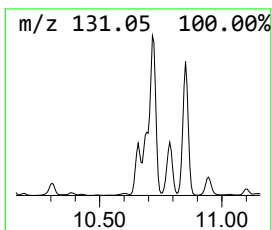
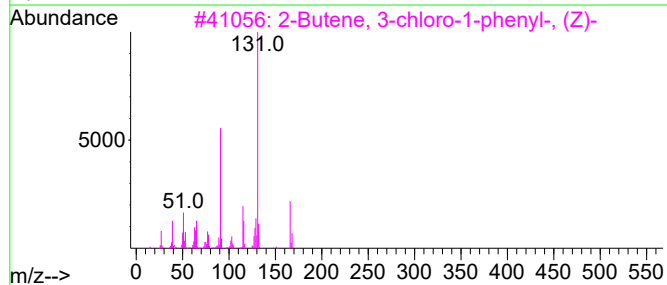
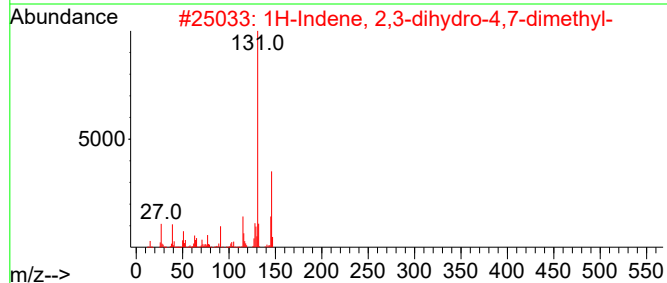
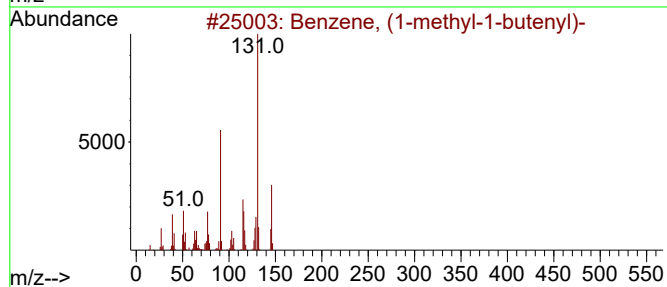
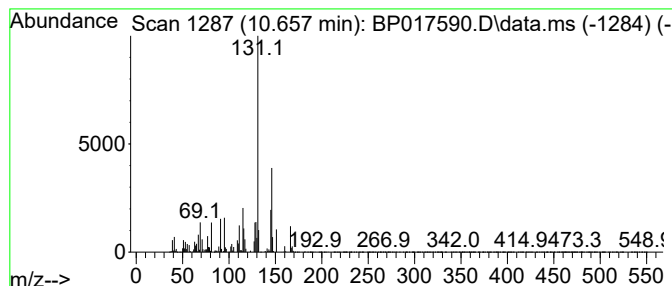
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 60 Benzene, (1-methyl-1-butenyl)- Concentration Rank 69

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.657	8.39 ng/ul	1650450	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	93
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3			2-Butene, 3-chloro-1-phenyl-, (Z)-	166	C10H11Cl	016608-68-7	93
4			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
5			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

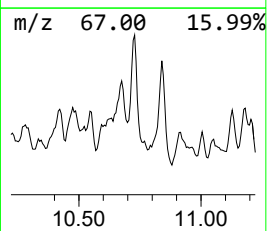
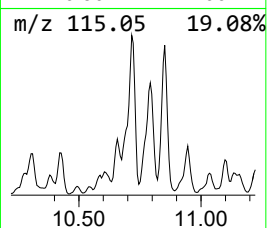
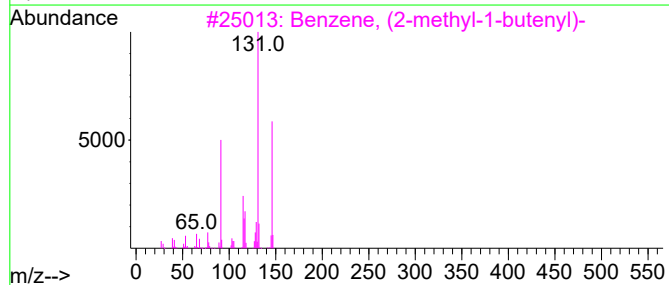
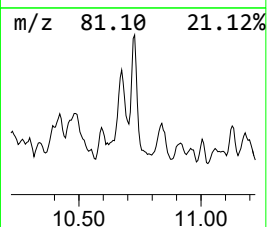
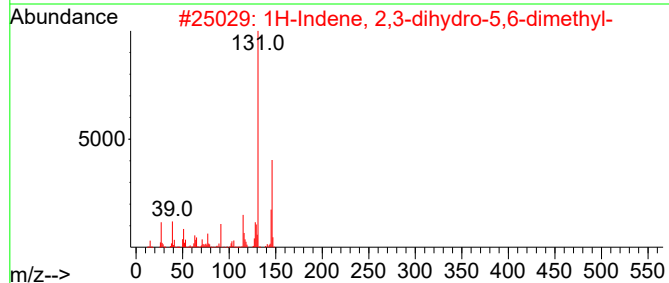
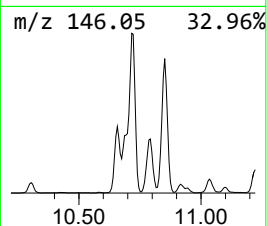
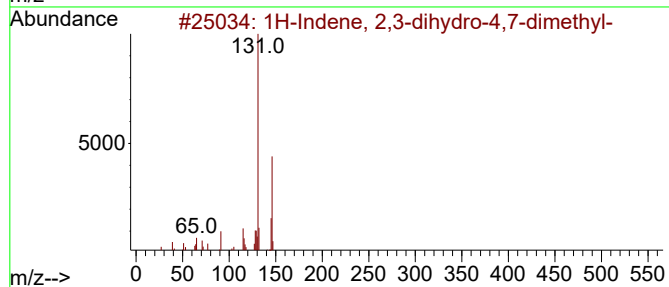
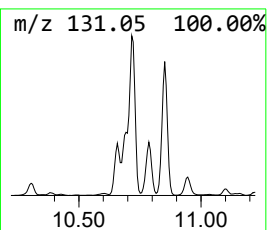
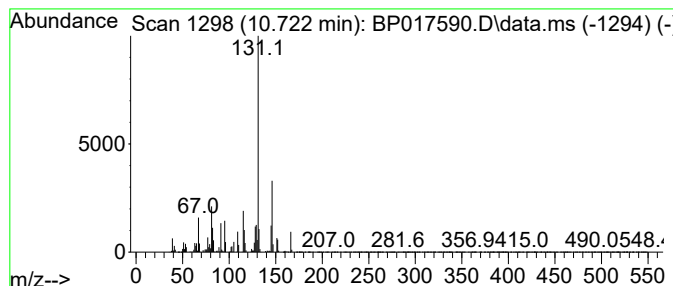
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 61 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.722	17.10 ng/ul	3364490	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
2			1H-Indene, 2,3-dihydro-5,6-dimet...	146	C11H14	001075-22-5	93
3			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	91
4			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	90
5			1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

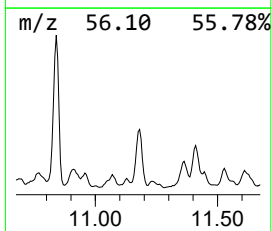
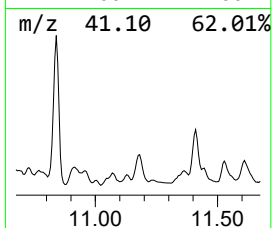
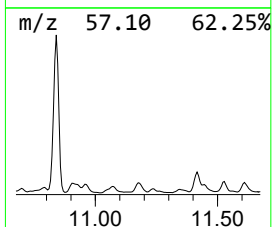
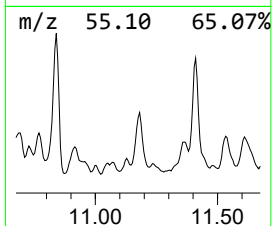
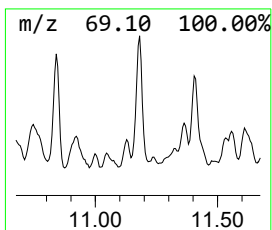
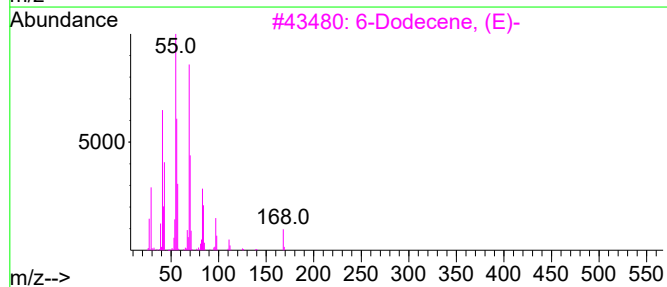
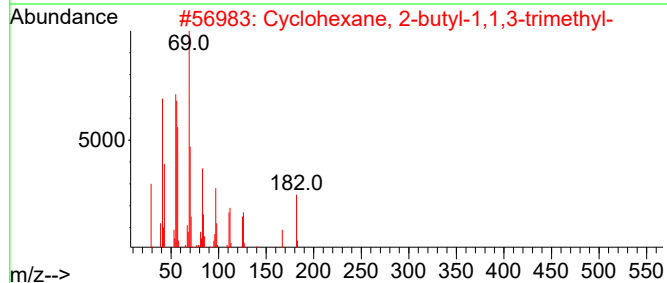
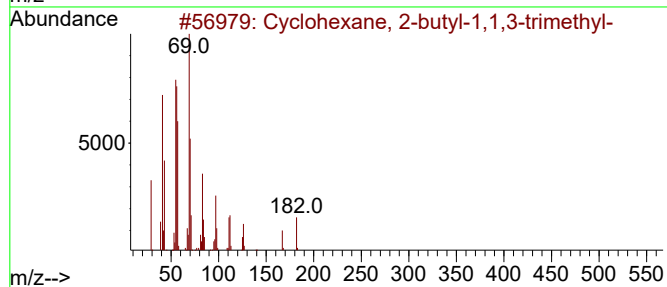
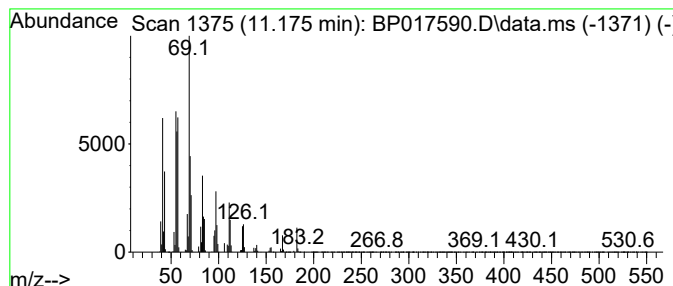
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 62 (DEL) Alkane: Cyclic11.175 Concentration Rank 73

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.175	8.17 ng/ul	1606780	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	96
2			Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	96
3			6-Dodecene, (E)-	168	C12H24	007206-17-9	76
4			Cyclopentane, 1,2-dibutyl-	182	C13H26	062199-52-4	72
5			Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

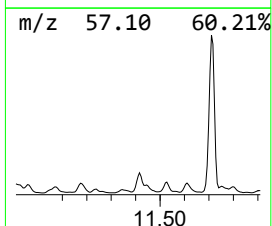
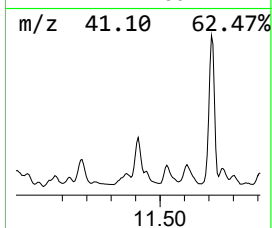
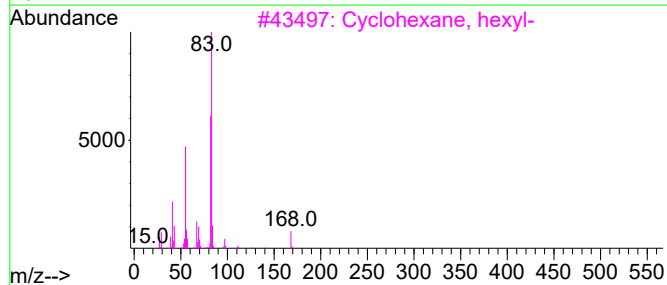
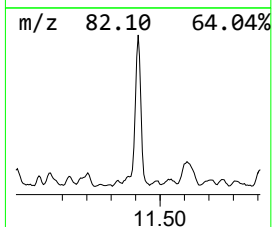
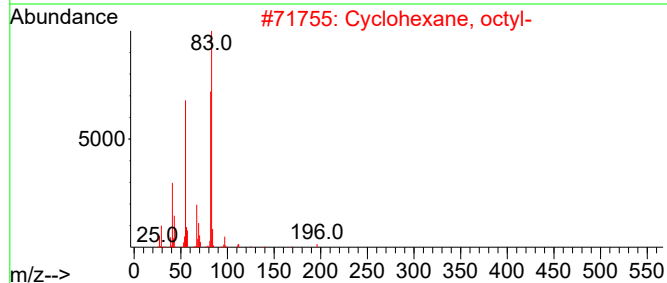
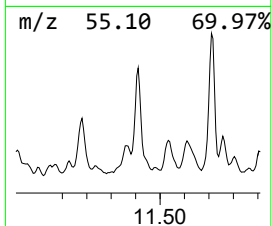
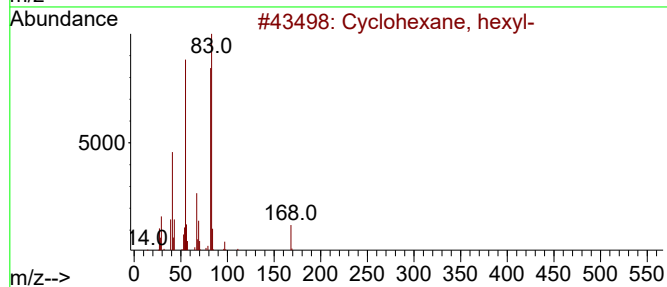
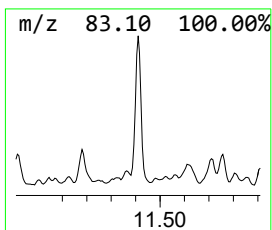
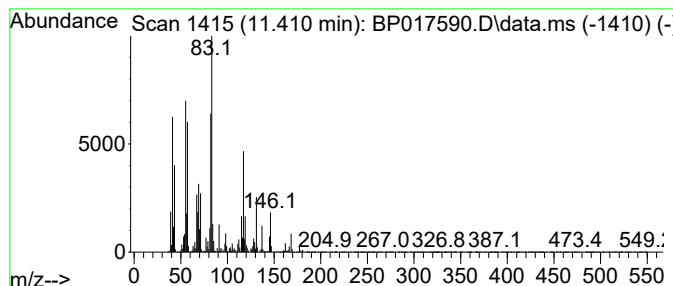
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 63 (DEL) Alkane: Cyclic11.410 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.410	18.12 ng/ul	3564210	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, hexyl-	168	C12H24	004292-75-5	55
2			Cyclohexane, octyl-	196	C14H28	001795-15-9	43
3			Cyclohexane, hexyl-	168	C12H24	004292-75-5	43
4			Cyclohexane, butyl-	140	C10H20	001678-93-9	43
5			Cyclohexane, ethyl-	112	C8H16	001678-91-7	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

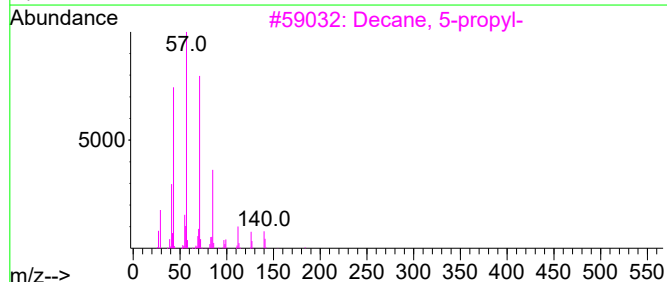
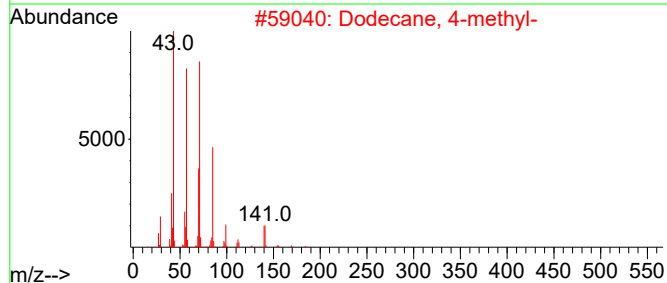
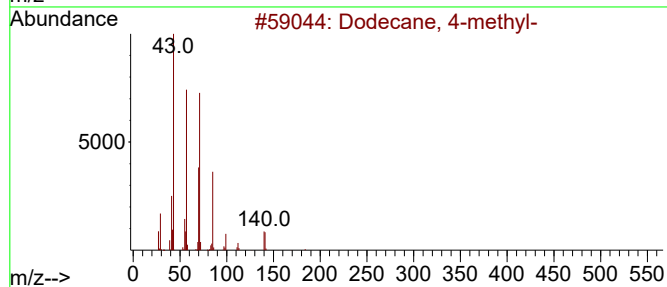
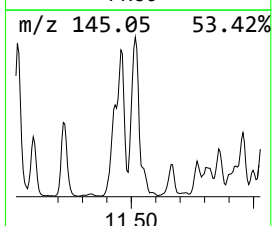
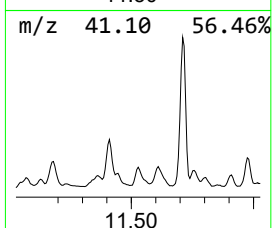
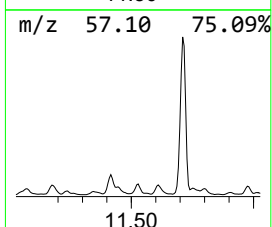
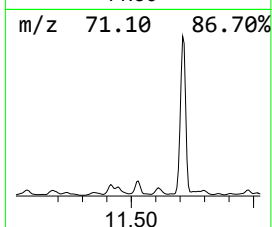
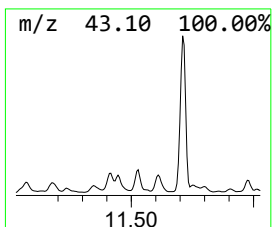
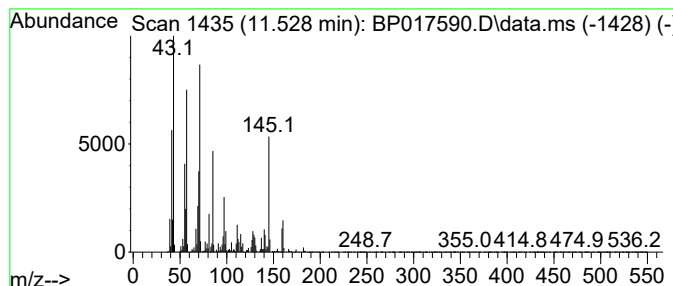
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 64 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.528	11.77 ng/ul	2316090	Naphthalene-d8	10.775	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 4-methyl-	184	C13H28	006117-97-1	91
2	Dodecane, 4-methyl-	184	C13H28	006117-97-1	90
3	Decane, 5-propyl-	184	C13H28	017312-62-8	49
4	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	38
5	Nonadecane, 9-methyl-	282	C20H42	013287-24-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

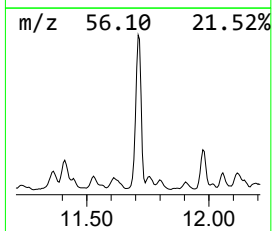
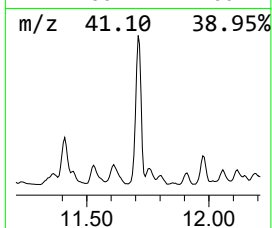
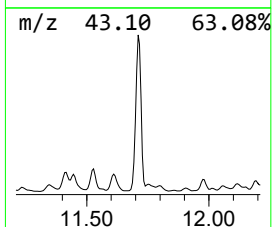
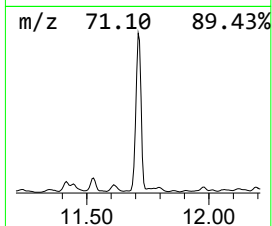
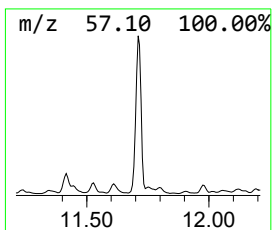
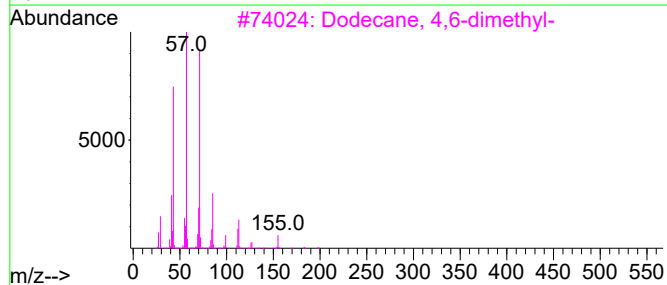
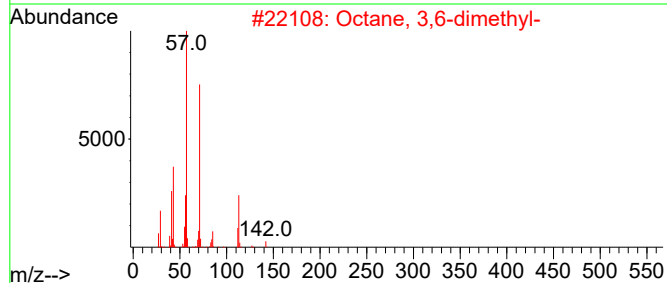
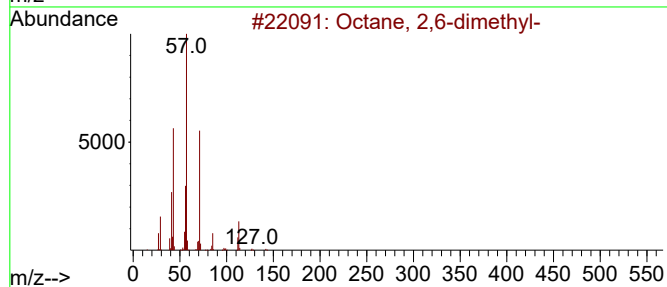
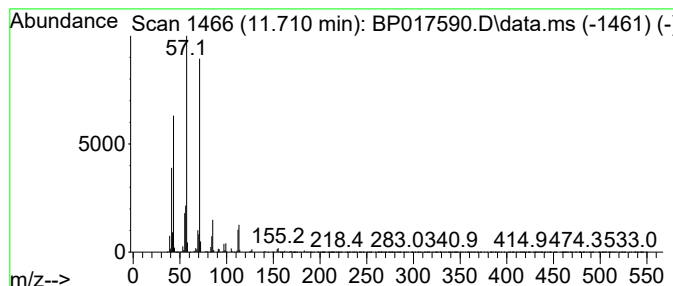
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 65 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.710	45.37 ng/ul	8925910	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 2,6-dimethyl-			142	C10H22	002051-30-1	90
2	Octane, 3,6-dimethyl-			142	C10H22	015869-94-0	83
3	Dodecane, 4,6-dimethyl-			198	C14H30	061141-72-8	64
4	Nonane, 2-methyl-5-propyl-			184	C13H28	031081-17-1	64
5	Heptane, 3-[(ethenyl)oxy)methyl]-			156	C10H20O	000103-44-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

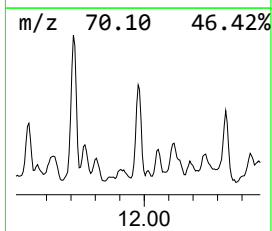
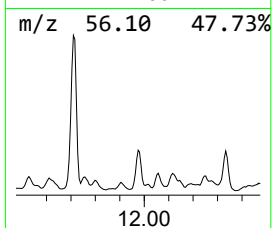
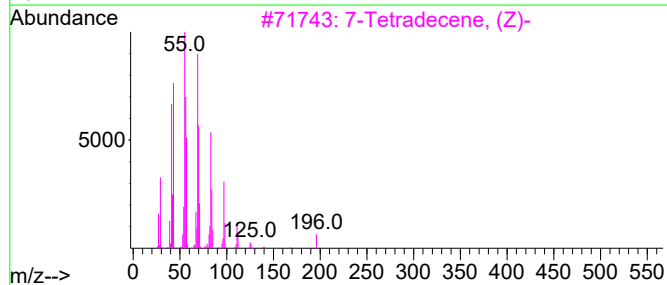
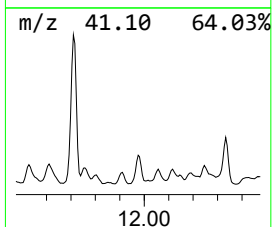
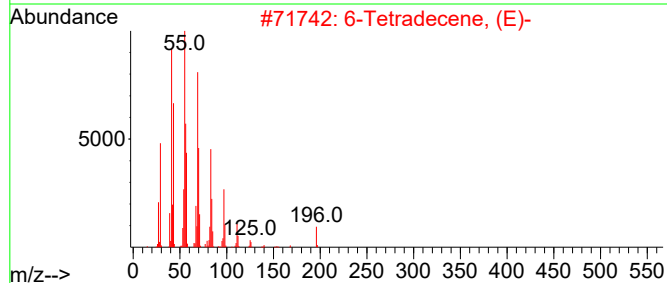
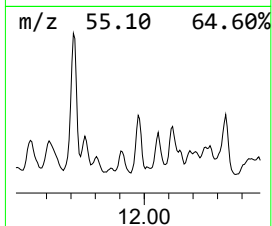
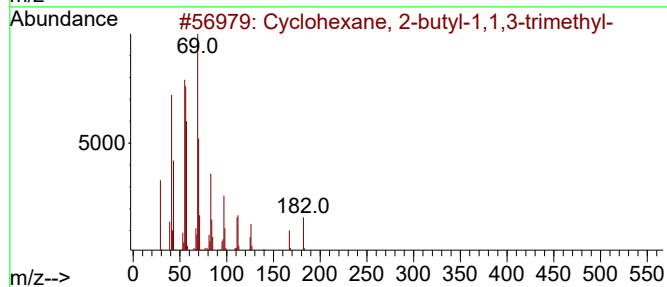
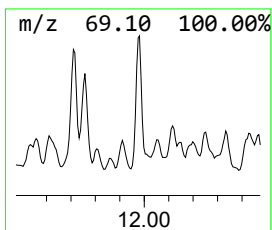
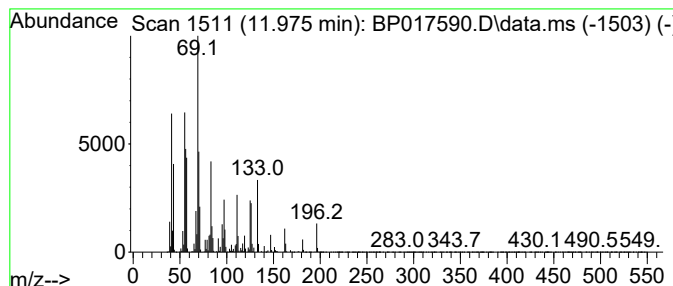
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 66 (DEL) Alkane: Cyclic11.975 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.975	14.85 ng/ul	2920590	Naphthalene-d8	10.775

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	81
2		6-Tetradecene, (E)-	196	C14H28	041446-64-4	76
3		7-Tetradecene, (Z)-	196	C14H28	041446-60-0	70
4		7-Tetradecene, (E)-	196	C14H28	041446-63-3	70
5		6-Tridecene, (E)-	182	C13H26	006434-76-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

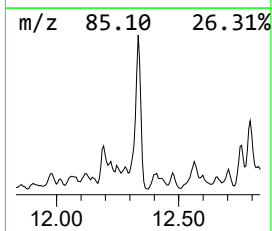
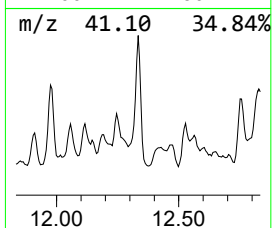
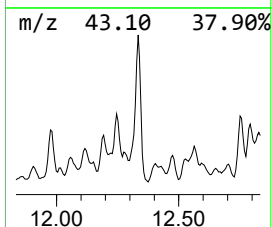
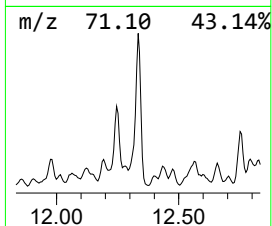
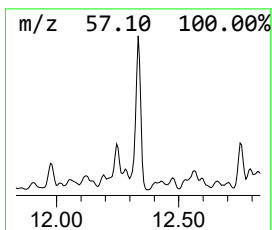
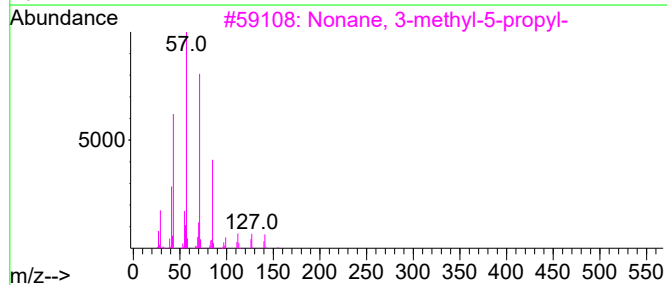
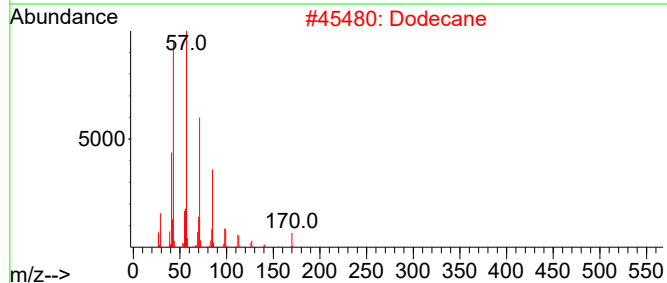
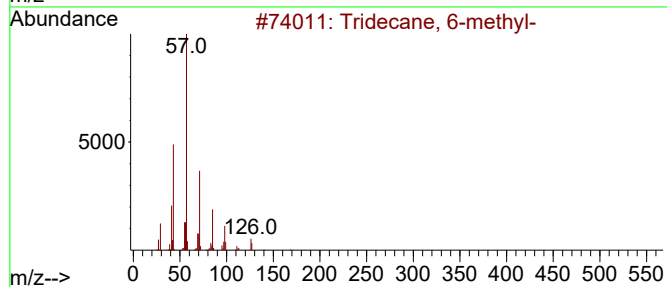
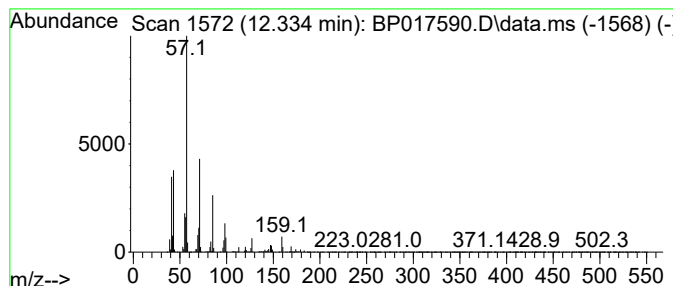
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 67 (DEL) Alkane: Straight-Chai... Concentration Rank 56

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.334	11.38 ng/ul	2238950	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 6-methyl-	198	C14H30	013287-21-3	90
2			Dodecane	170	C12H26	000112-40-3	59
3			Nonane, 3-methyl-5-propyl-	184	C13H28	031081-18-2	59
4			Undecane, 5-ethyl-	184	C13H28	017453-94-0	59
5			Pentadecane	212	C15H32	000629-62-9	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

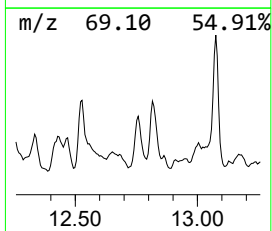
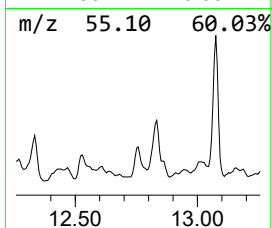
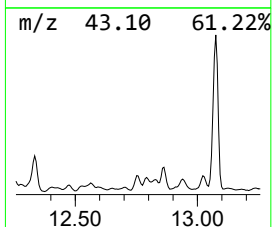
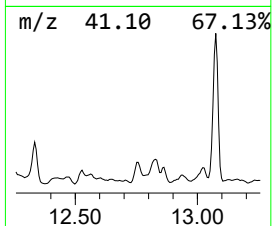
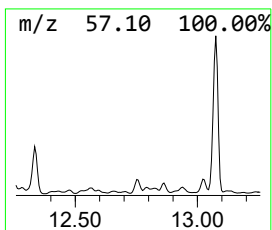
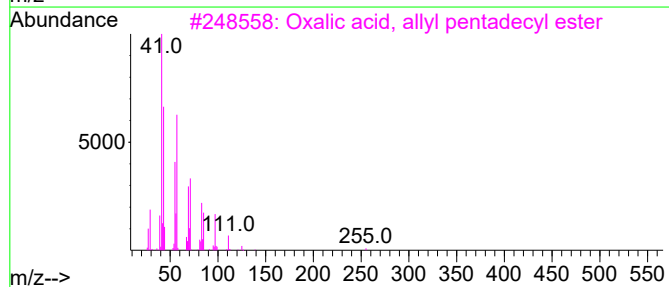
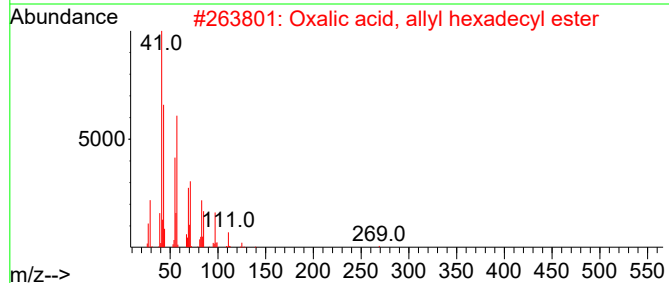
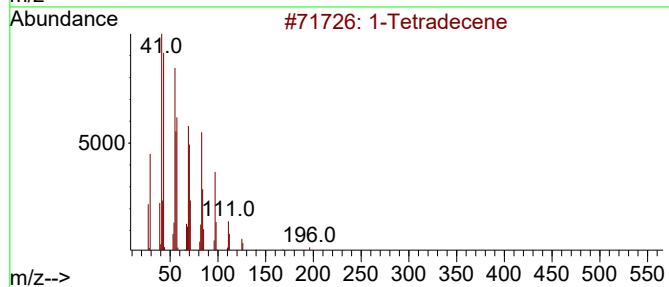
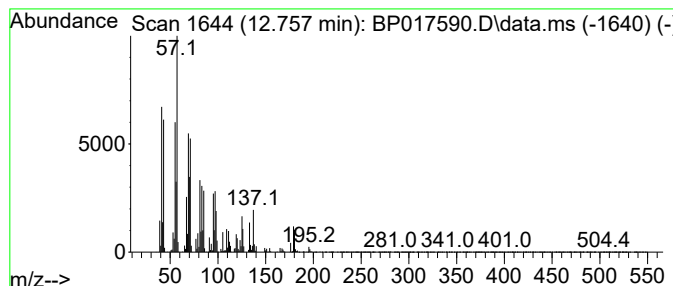
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 68 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.757	14.48 ng/ul	2010900	Acenaphthene-d10		14.628	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Tetradecene		196	C14H28	001120-36-1	44
2	Oxalic acid, allyl hexadecyl ester		354	C21H38O4	1000309-24-4	43
3	Oxalic acid, allyl pentadecyl ester		340	C20H36O4	1000309-24-3	38
4	Decane, 2,3,8-trimethyl-		184	C13H28	062238-14-6	22
5	Undecane		156	C11H24	001120-21-4	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

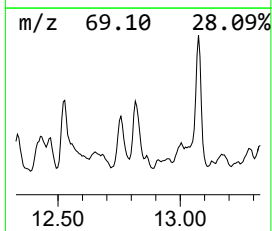
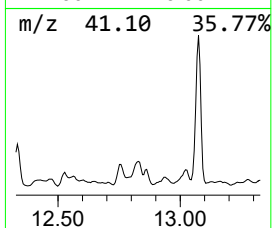
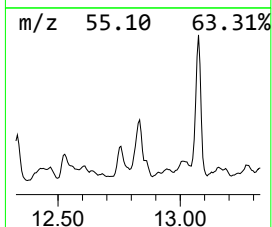
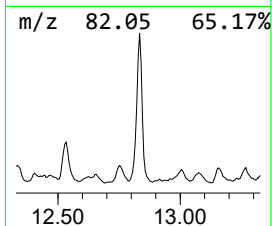
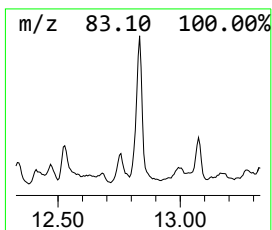
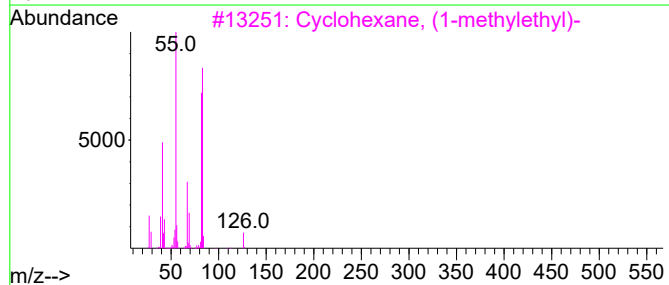
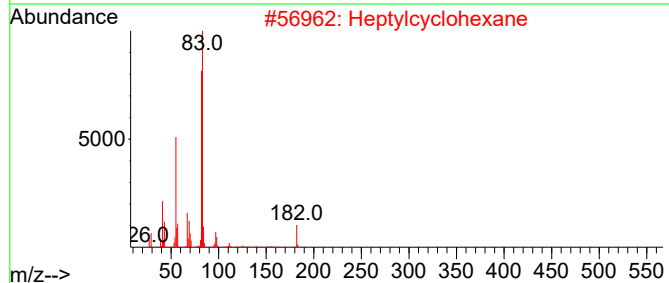
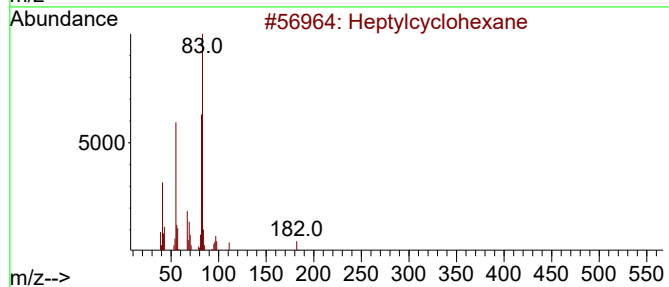
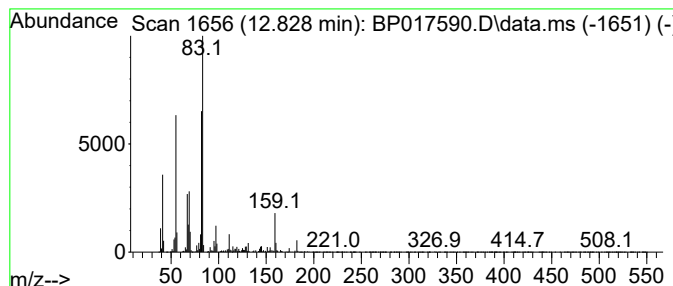
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 69 (DEL) Alkane: Cyclic12.828 Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.828	15.88 ng/ul	2205300	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptylcyclohexane	182	C13H26	005617-41-4	80
2			Heptylcyclohexane	182	C13H26	005617-41-4	64
3			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	59
5			Cyclohexane, pentyl-	154	C11H22	004292-92-6	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

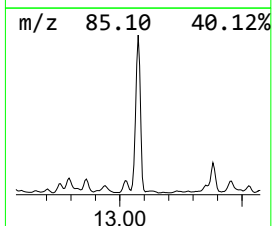
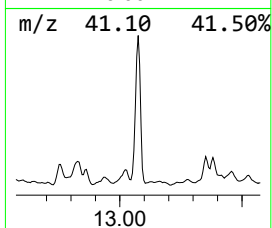
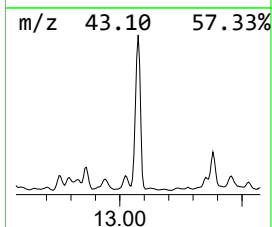
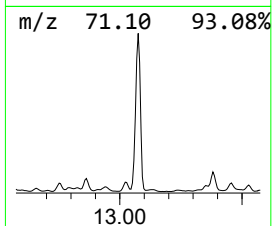
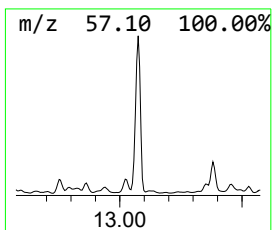
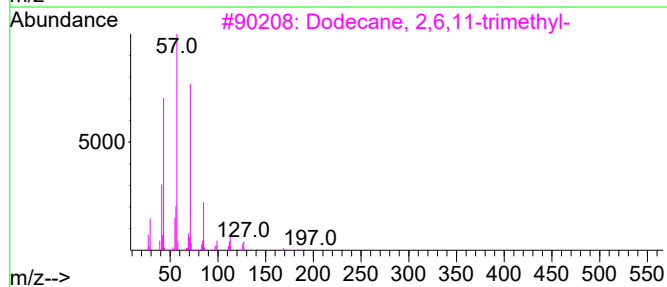
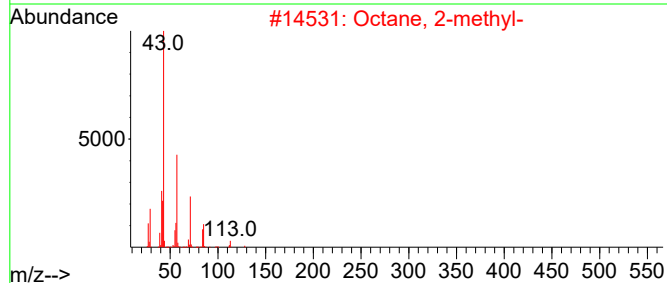
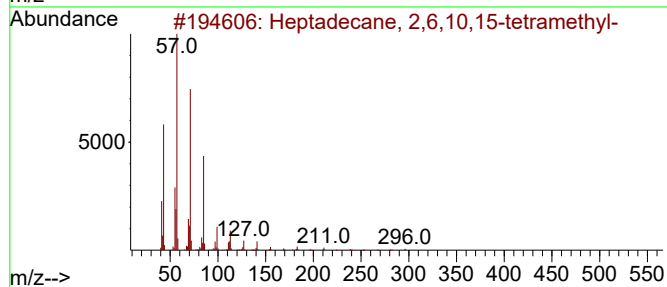
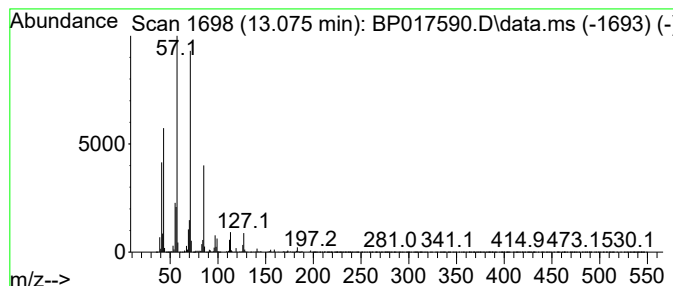
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 70 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.075	73.01 ng/ul	10136100	Acenaphthene-d10	14.628

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
2		Octane, 2-methyl-	128	C9H20	003221-61-2	87
3		Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	80
4		Dodecane, 2,7,10-trimethyl-	212	C15H32	074645-98-0	80
5		Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

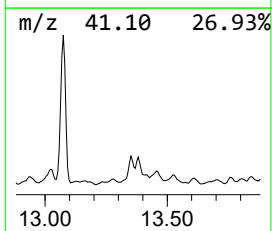
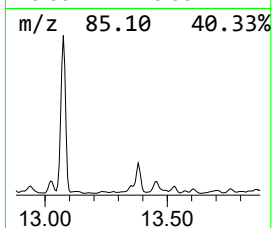
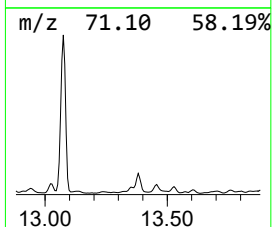
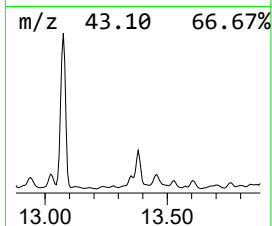
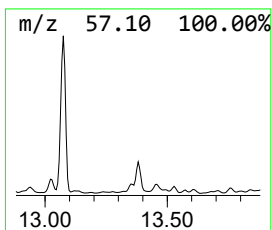
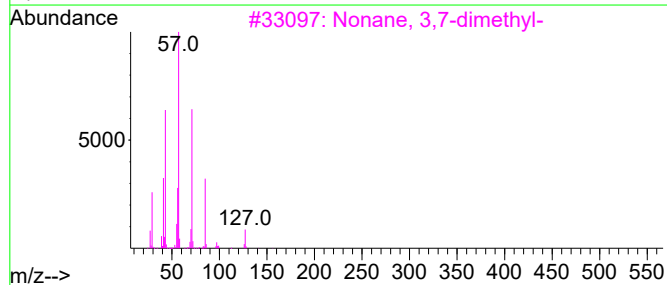
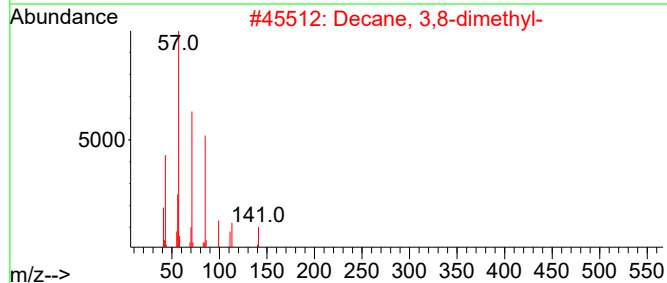
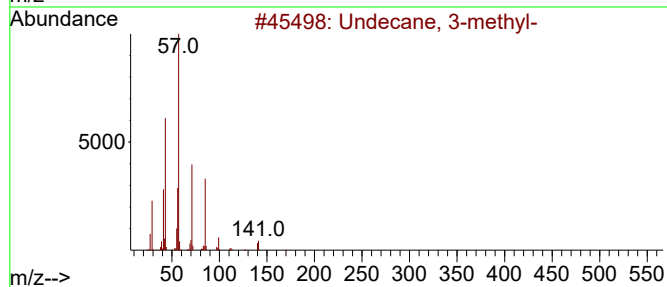
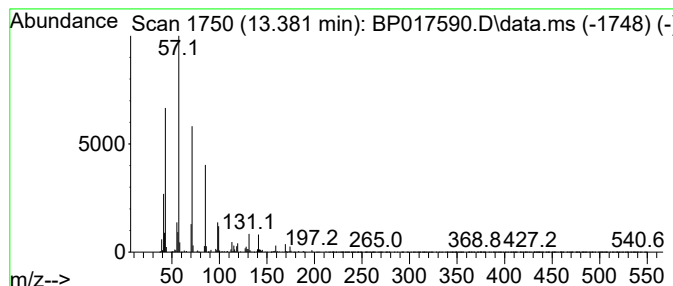
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 71 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.381	12.16 ng/ul	1688670	Acenaphthene-d10	14.628	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 3-methyl-	170	C12H26	001002-43-3	64
2	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	62
3	Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	58
4	Decane, 1-iodo-	268	C10H21I	002050-77-3	53
5	1,3-Propanediol, decyl ethyl ether	244	C15H32O2	1000406-35-0	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

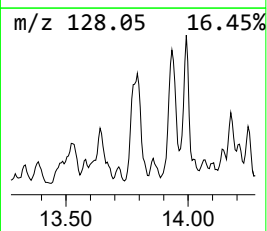
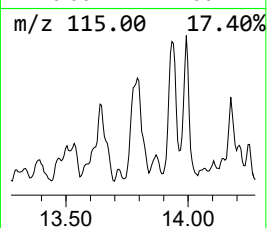
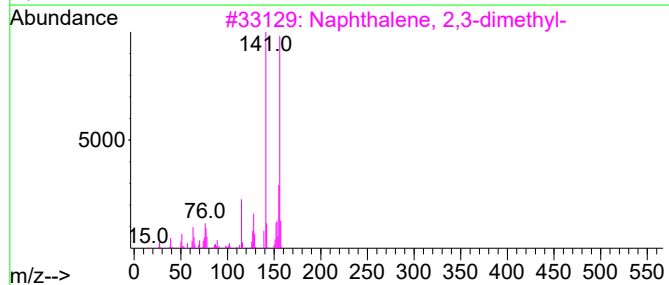
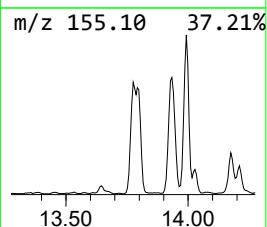
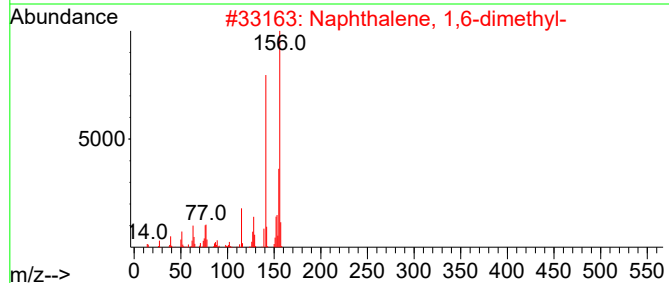
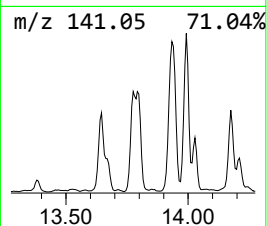
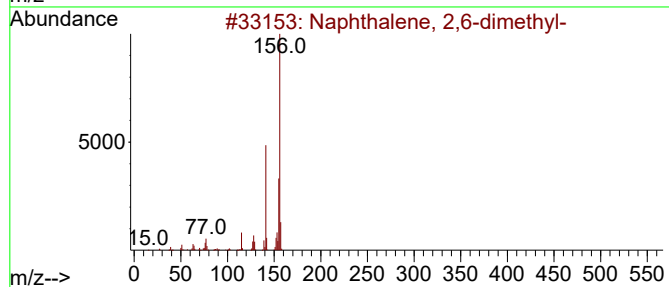
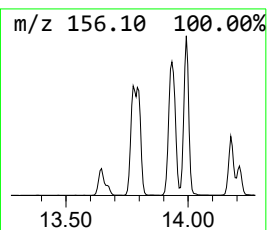
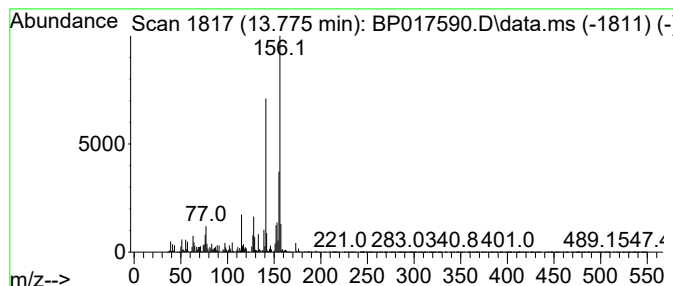
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 72 Naphthalene, 2,6-dimethyl- Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.775	14.46 ng/ul	2007630	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

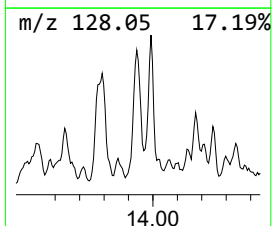
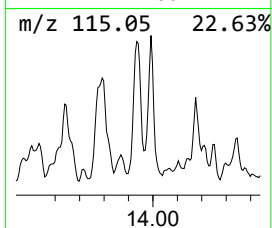
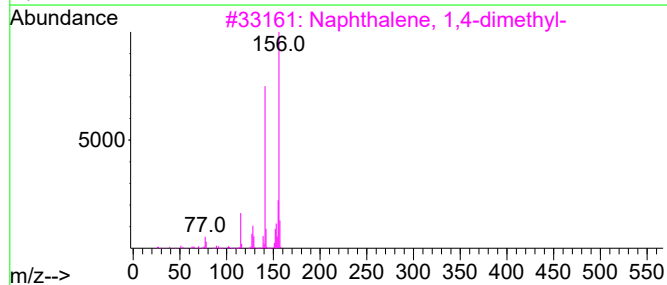
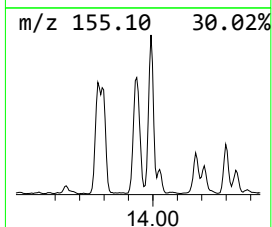
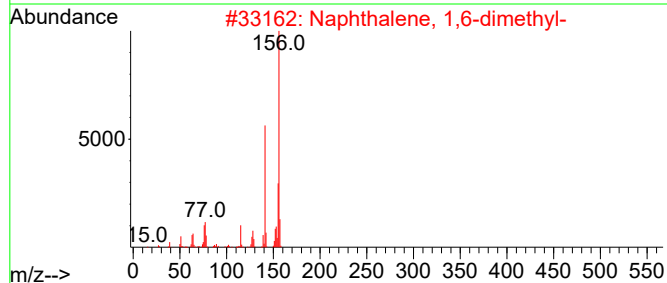
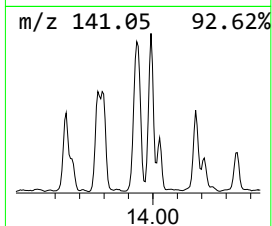
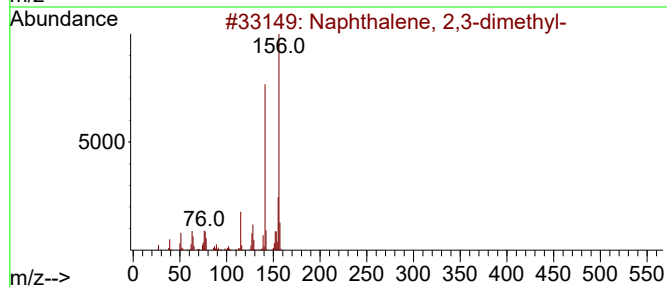
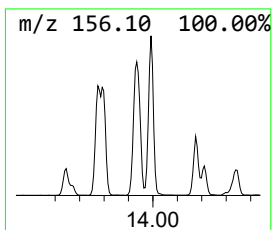
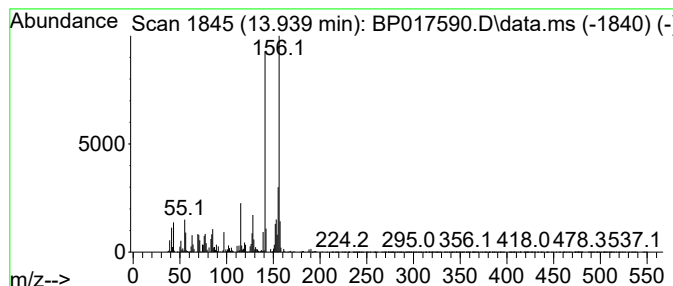
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 73 Naphthalene, 2,3-dimethyl- Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.940	17.41 ng/ul	2417520	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

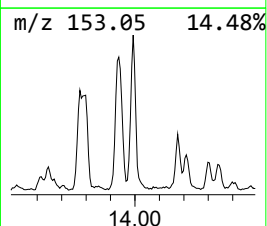
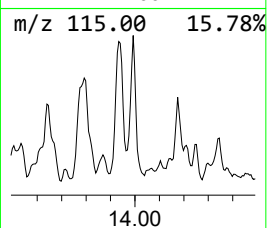
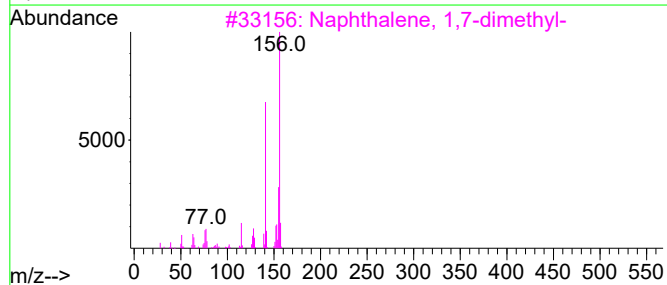
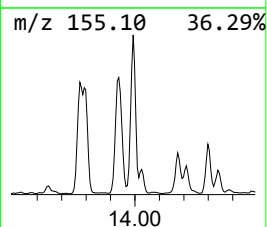
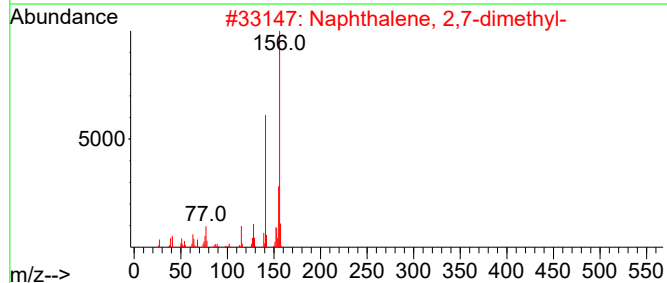
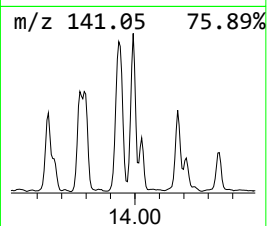
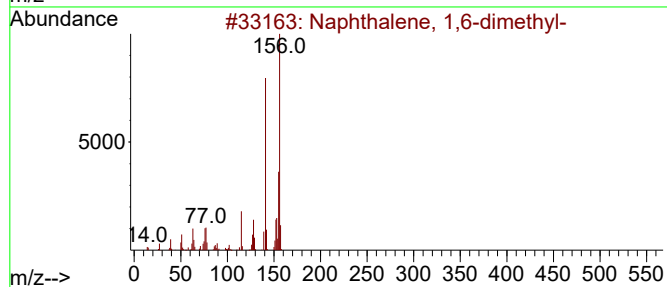
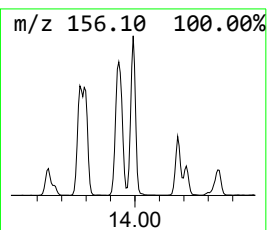
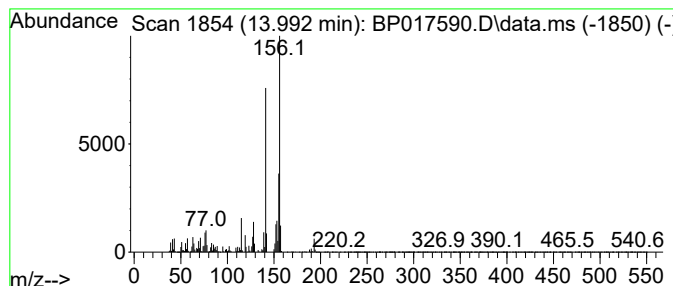
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 74 Naphthalene, 1,6-dimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.992	17.35 ng/ul	2408650	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
3			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
5			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

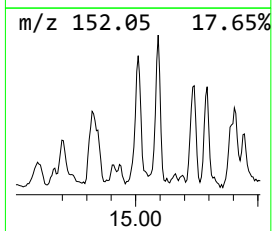
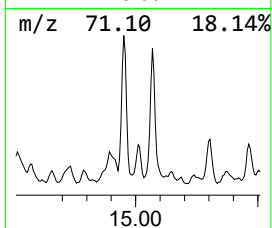
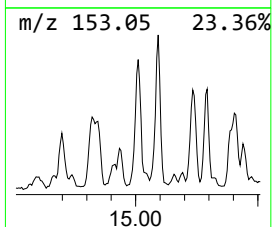
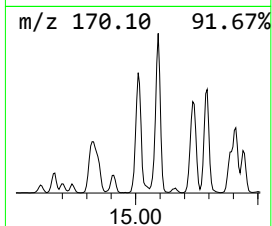
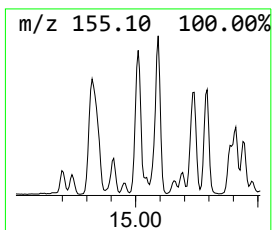
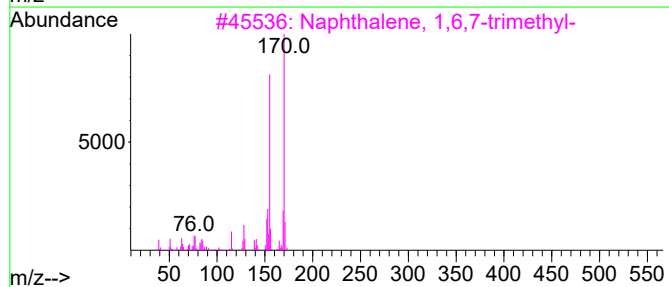
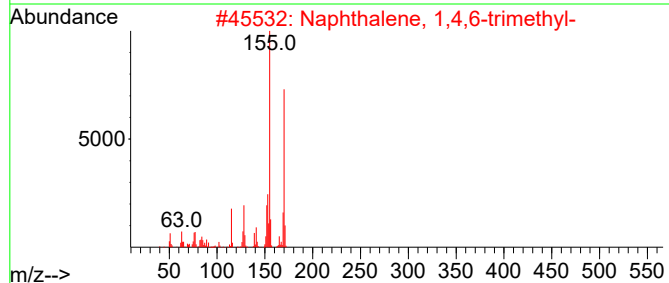
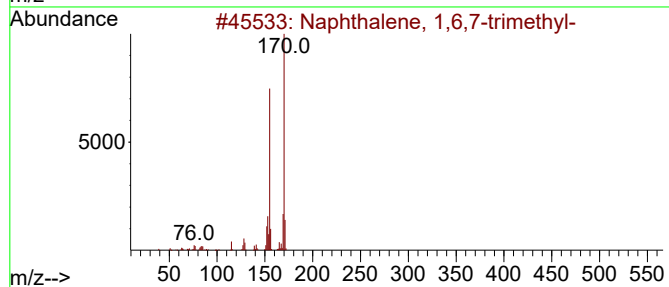
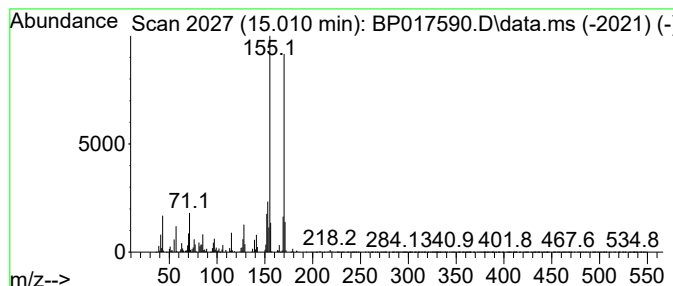
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 75 Naphthalene, 1,6,7-trimethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.010	13.52 ng/ul	1876720	Acenaphthene-d10	14.628	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96
2	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
5	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

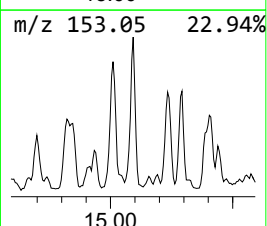
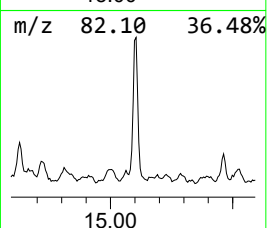
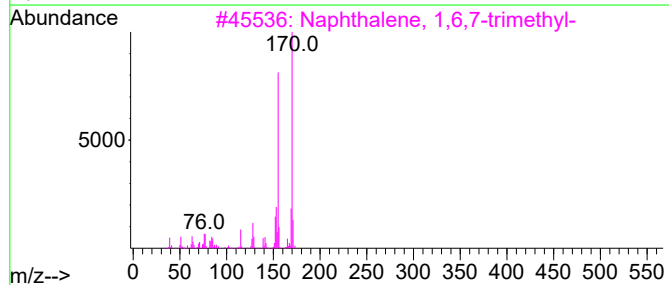
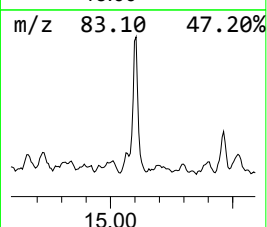
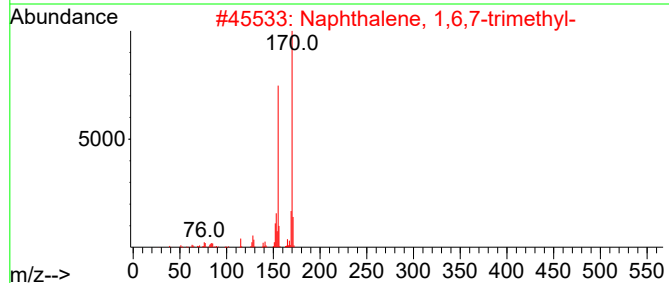
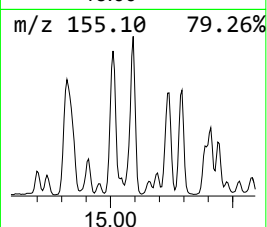
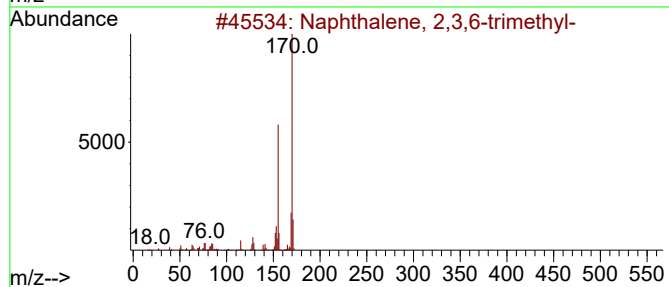
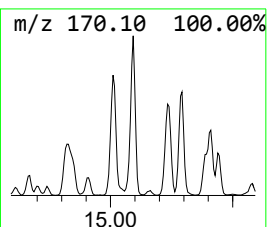
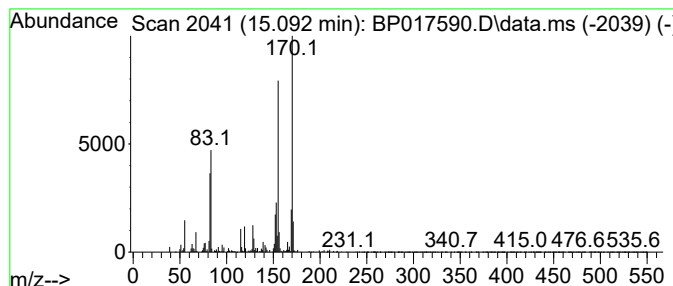
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 76 Naphthalene, 2,3,6-trimethyl- Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.092	11.46 ng/ul	1590690	Acenaphthene-d10	14.628	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
3	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	90
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
5	Naphthalene, 1,4,5-trimethyl-	170	C13H14	002131-41-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

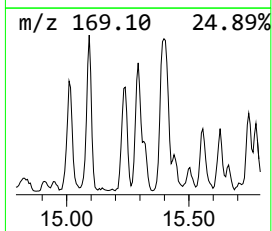
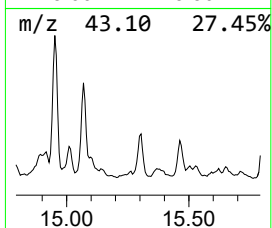
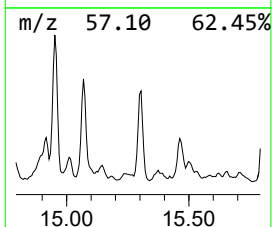
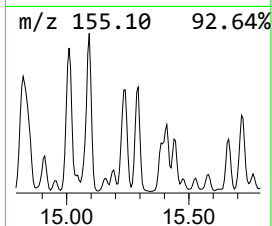
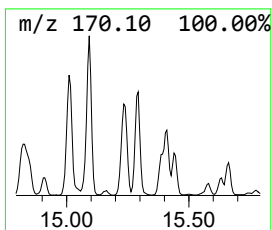
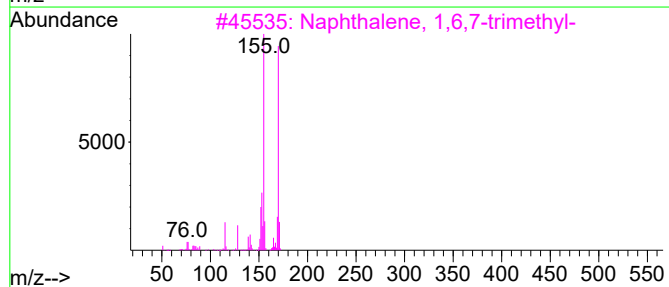
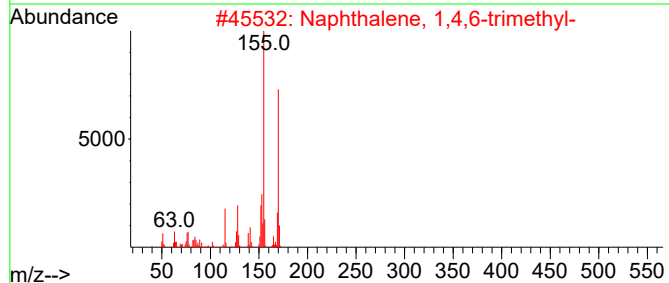
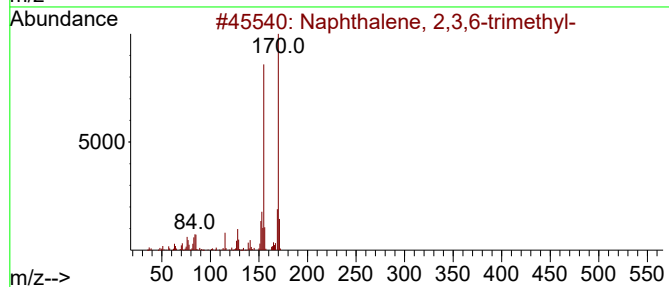
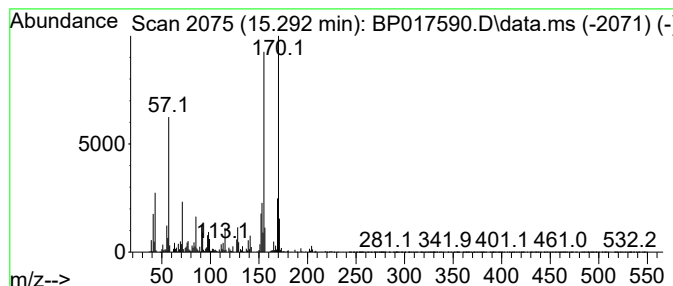
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 77 Naphthalene, 1,4,6-trimethyl- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.292	13.15 ng/ul	1825680	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	96
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

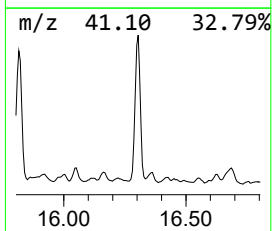
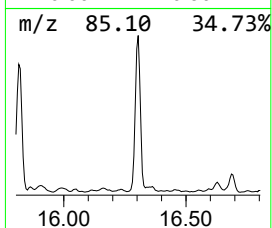
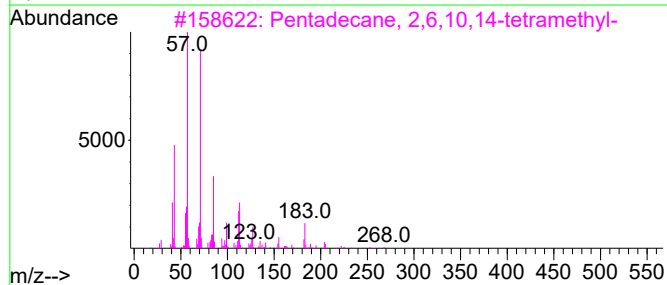
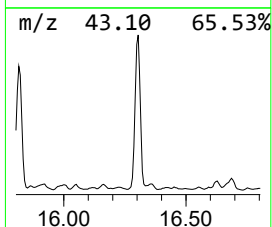
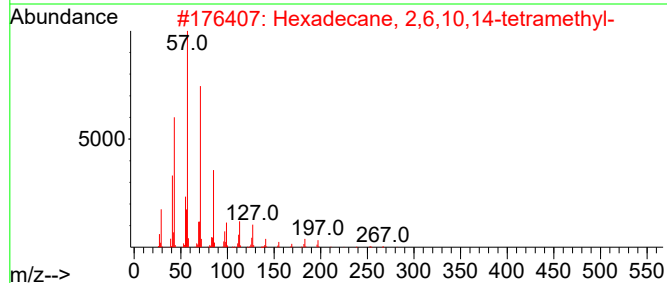
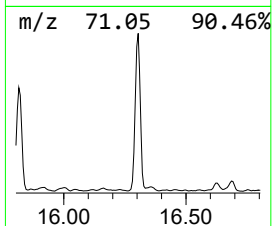
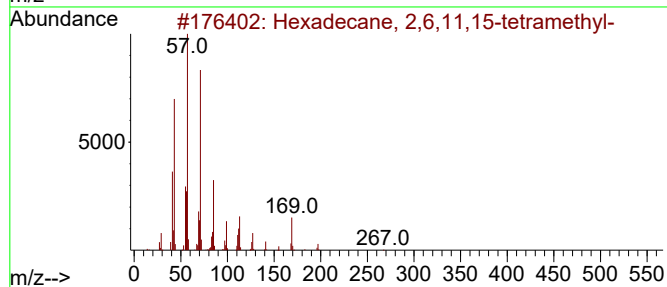
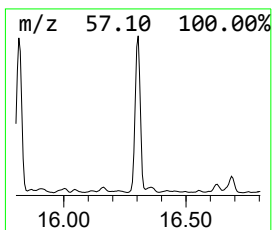
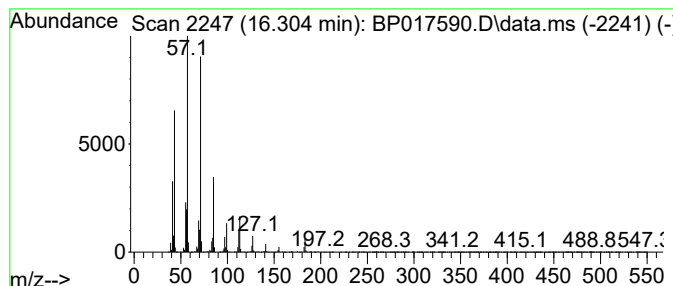
Instrument :
BNA_P
ClientSampleId :
BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 78 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
16.304	97.07 ng/ul	7839360	Phenanthrene-d10	17.416	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,11,15-tetramethyl-	282	C20H42	000504-44-9	91
2	Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	90
3	Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	90
4	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	72
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

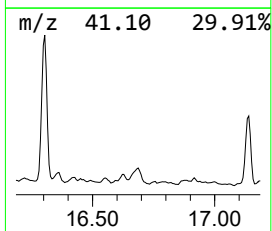
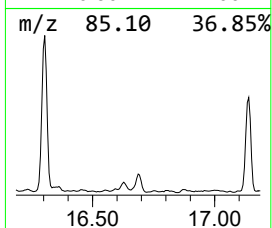
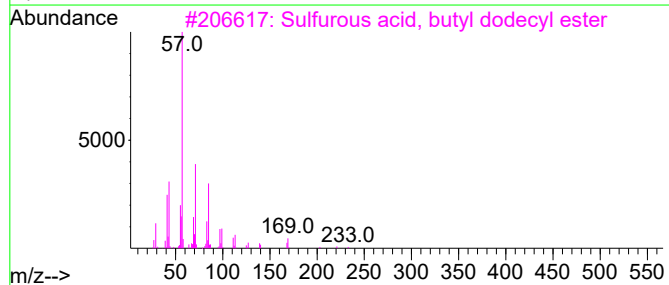
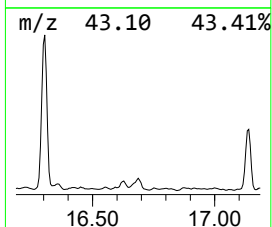
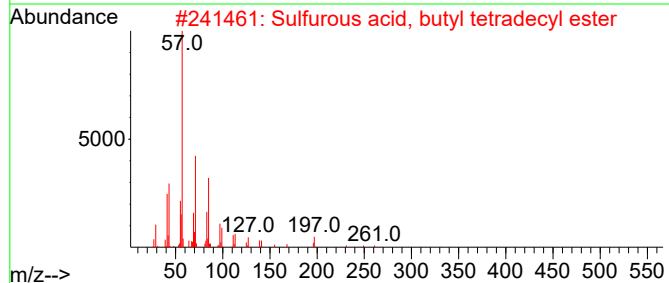
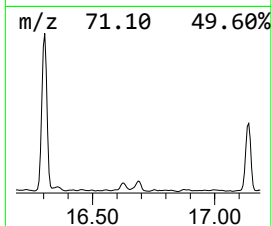
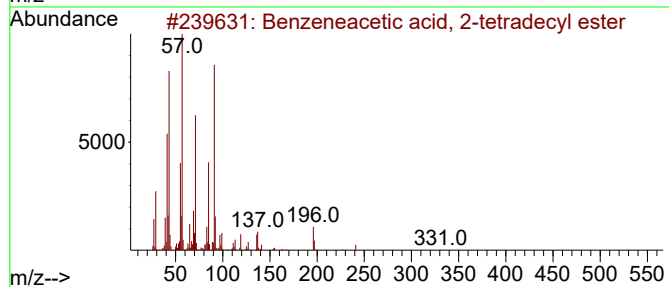
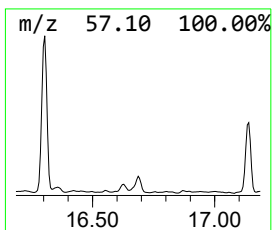
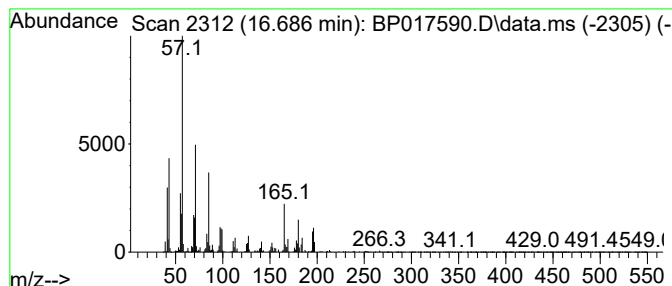
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 79 Benzeneacetic acid, 2-tetra... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.686	20.87 ng/ul	1685330	Phenanthrene-d10	17.416

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzeneacetic acid, 2-tetradecyl...	332	C22H36O2	1000282-02-4	53
2			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	49
3			Sulfurous acid, butyl dodecyl ester	306	C16H34O3S	1000309-17-9	49
4			Tetratetracontane	619	C44H90	007098-22-8	46
5			Hexacosane	366	C26H54	000630-01-3	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
Data File : BP017590.D
Acq On : 06 Oct 2023 06:43
Operator : MA/JU
Sample : 04588-09
Misc :
ALS Vial : 67 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BBHT7

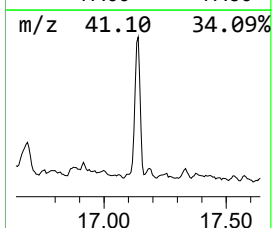
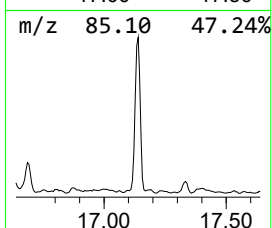
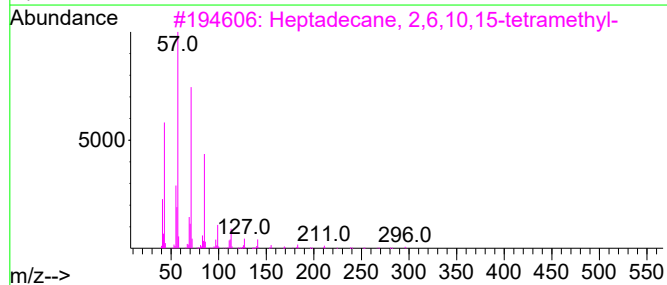
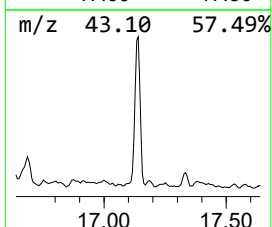
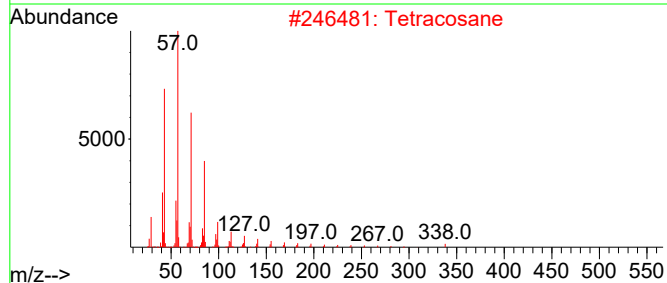
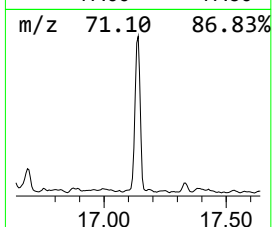
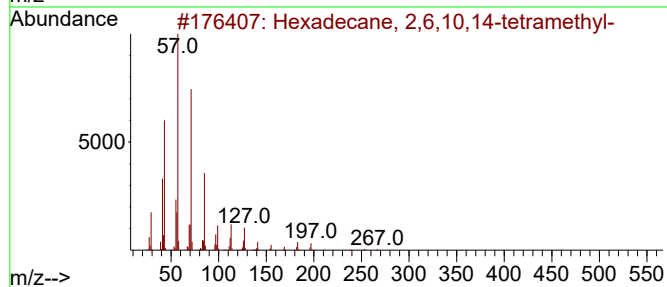
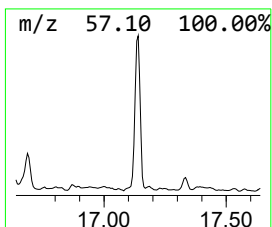
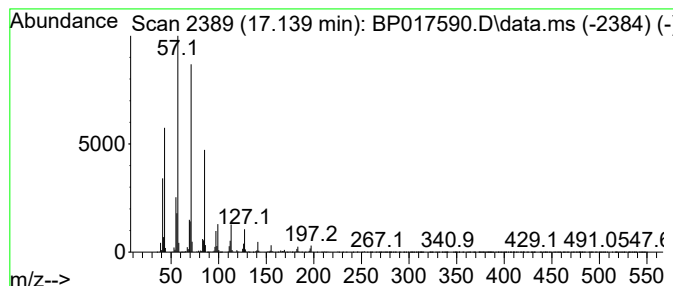
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST20.L
TIC Integration Parameters: LSCINT.P

Peak Number 80 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.139	43.27 ng/ul	3494790	Phenanthrene-d10	17.416

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Tetracosane	338	C24H50	000646-31-1	91
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
4		Hexadecane, 1-iodo-	352	C16H33I	000544-77-4	87
5		Octadecane, 1-iodo-	380	C18H37I	000629-93-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100423\
 Data File : BP017590.D
 Acq On : 06 Oct 2023 06:43
 Operator : MA/JU
 Sample : 04588-09
 Misc :
 ALS Vial : 67 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BBHT7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100423.MA.M
 Quant Title : SVOA CALIBRATION

TIC Library : C:\DATABASE\NIST0.L
 TIC Integration Parameters: LSCINT.P

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	5.263	8.7	ng/ul	2379360	1	7.940	5448060	20.0
(DEL) Alkane: S...	5.399	8.9	ng/ul	2429420	1	7.940	5448060	20.0
(DEL) Alkane: C...	5.705	5.8	ng/ul	1578760	1	7.940	5448060	20.0
(DEL) Alkane: C...	5.799	13.1	ng/ul	3557810	1	7.940	5448060	20.0
(DEL) Alkane: C...	5.863	6.6	ng/ul	1790720	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.093	14.7	ng/ul	3992870	1	7.940	5448060	20.0
Sulfurous acid,...	6.122	10.7	ng/ul	2924130	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.222	14.0	ng/ul	3810350	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.310	9.5	ng/ul	2596680	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.393	48.7	ng/ul	13276900	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.440	9.3	ng/ul	2546280	1	7.940	5448060	20.0
Ethanone, 1-cyc...	6.493	53.3	ng/ul	14517000	1	7.940	5448060	20.0
(DEL) Alkane: C...	6.599	7.7	ng/ul	2097140	1	7.940	5448060	20.0
Cyclohexane, 1,...	6.669	10.4	ng/ul	2837250	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.710	8.2	ng/ul	2231510	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.734	13.2	ng/ul	3588900	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.793	7.8	ng/ul	2115480	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.834	30.8	ng/ul	8388620	1	7.940	5448060	20.0
(DEL) Alkane: S...	6.887	9.0	ng/ul	2462560	1	7.940	5448060	20.0
(DEL) Alkane: C...	6.940	8.2	ng/ul	2243060	1	7.940	5448060	20.0
(DEL) Alkane: C...	7.005	52.3	ng/ul	14256000	1	7.940	5448060	20.0
1,5-Hexadiene, ...	7.146	6.0	ng/ul	1626940	1	7.940	5448060	20.0
(DEL) Alkane: C...	7.175	15.0	ng/ul	4093370	1	7.940	5448060	20.0
(DEL) Alkane: C...	7.346	30.3	ng/ul	8249760	1	7.940	5448060	20.0
1-Propanone, 1-...	7.387	12.1	ng/ul	3305670	1	7.940	5448060	20.0
Mesitylene	7.546	10.4	ng/ul	2841390	1	7.940	5448060	20.0
(DEL) Alkane: C...	7.587	11.9	ng/ul	3251950	1	7.940	5448060	20.0
Sulfurous acid,...	7.640	8.5	ng/ul	2307220	1	7.940	5448060	20.0
Cyclobutylcarbo...	7.675	9.7	ng/ul	2632230	1	7.940	5448060	20.0
unknown-01	7.769	25.8	ng/ul	7016990	1	7.940	5448060	20.0
(DEL) Alkane: S...	7.834	64.8	ng/ul	17666500	1	7.940	5448060	20.0
unknown-02	8.046	29.7	ng/ul	8080700	1	7.940	5448060	20.0
(DEL) Alkane: S...	8.110	11.9	ng/ul	3242890	1	7.940	5448060	20.0
(DEL) Alkane: C...	8.169	35.1	ng/ul	9552070	1	7.940	5448060	20.0
(DEL) Alkane: C...	8.216	13.4	ng/ul	3650610	1	7.940	5448060	20.0
(DEL) Alkane: C...	8.310	18.8	ng/ul	5125280	1	7.940	5448060	20.0
(DEL) Alkane: S...	8.387	39.5	ng/ul	10773900	1	7.940	5448060	20.0
(DEL) Alkane: S...	8.446	33.1	ng/ul	9018730	1	7.940	5448060	20.0
(DEL) Alkane: S...	8.522	14.3	ng/ul	3894810	1	7.940	5448060	20.0
Benzene, (2-met...	8.552	11.5	ng/ul	3139650	1	7.940	5448060	20.0
Naphthalene, de...	8.734	36.3	ng/ul	9891930	1	7.940	5448060	20.0
Benzene, 4-ethy...	8.863	7.0	ng/ul	1896420	1	7.940	5448060	20.0
(DEL) Alkane: S...	8.957	15.1	ng/ul	4123380	1	7.940	5448060	20.0
Ethanone, 1-(1-...	8.993	24.0	ng/ul	6528970	1	7.940	5448060	20.0
Benzene, 1-ethy...	9.022	30.6	ng/ul	8346760	1	7.940	5448060	20.0
(DEL) Alkane: S...	9.081	12.3	ng/ul	3339150	1	7.940	5448060	20.0
(DEL) Alkane: C...	9.204	8.2	ng/ul	2240530	1	7.940	5448060	20.0
Benzene, 1-ethy...	9.246	13.8	ng/ul	3746630	1	7.940	5448060	20.0
(DEL) Alkane: S...	9.346	20.1	ng/ul	5471930	1	7.940	5448060	20.0
(DEL) Alkane: S...	9.510	26.6	ng/ul	5231370	2	10.775	3934480	20.0
Benzene, 1,2,4,...	9.551	19.1	ng/ul	3762490	2	10.775	3934480	20.0
trans-Decalin, ...	9.628	46.7	ng/ul	9188160	2	10.775	3934480	20.0

(DEL) Alkane: S...	9.746	10.4	ng/ul	2055880	2	10.775	3934480	20.0
unknown-03	9.922	36.1	ng/ul	7101430	2	10.775	3934480	20.0
(DEL) Alkane: S...	10.016	22.4	ng/ul	4404490	2	10.775	3934480	20.0
3-Phenylbut-1-ene	10.134	50.9	ng/ul	10005300	2	10.775	3934480	20.0
Ethanone, 1-(2,...	10.193	15.6	ng/ul	3076460	2	10.775	3934480	20.0
(DEL) Alkane: C...	10.593	8.6	ng/ul	1687450	2	10.775	3934480	20.0
Benzene, (1-met...	10.657	8.4	ng/ul	1650450	2	10.775	3934480	20.0
1H-Indene, 2,3-...	10.722	17.1	ng/ul	3364490	2	10.775	3934480	20.0
(DEL) Alkane: C...	11.175	8.2	ng/ul	1606780	2	10.775	3934480	20.0
(DEL) Alkane: C...	11.410	18.1	ng/ul	3564210	2	10.775	3934480	20.0
(DEL) Alkane: S...	11.528	11.8	ng/ul	2316090	2	10.775	3934480	20.0
(DEL) Alkane: S...	11.710	45.4	ng/ul	8925910	2	10.775	3934480	20.0
(DEL) Alkane: C...	11.975	14.9	ng/ul	2920590	2	10.775	3934480	20.0
(DEL) Alkane: S...	12.334	11.4	ng/ul	2238950	2	10.775	3934480	20.0
(DEL) Alkane: S...	12.757	14.5	ng/ul	2010900	3	14.628	2776800	20.0
(DEL) Alkane: C...	12.828	15.9	ng/ul	2205300	3	14.628	2776800	20.0
(DEL) Alkane: S...	13.075	73.0	ng/ul	10136100	3	14.628	2776800	20.0
(DEL) Alkane: S...	13.381	12.2	ng/ul	1688670	3	14.628	2776800	20.0
Naphthalene, 2,...	13.775	14.5	ng/ul	2007630	3	14.628	2776800	20.0
Naphthalene, 2,...	13.940	17.4	ng/ul	2417520	3	14.628	2776800	20.0
Naphthalene, 1,...	13.992	17.4	ng/ul	2408650	3	14.628	2776800	20.0
Naphthalene, 1,...	15.010	13.5	ng/ul	1876720	3	14.628	2776800	20.0
Naphthalene, 2,...	15.092	11.5	ng/ul	1590690	3	14.628	2776800	20.0
Naphthalene, 1,...	15.292	13.2	ng/ul	1825680	3	14.628	2776800	20.0
(DEL) Alkane: S...	16.304	97.1	ng/ul	7839360	4	17.416	1615200	20.0
Benzeneacetic a...	16.686	20.9	ng/ul	1685330	4	17.416	1615200	20.0
(DEL) Alkane: S...	17.139	43.3	ng/ul	3494790	4	17.416	1615200	20.0

Instrument :

BNA_P

ClientSampleId :

BBHT7