

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
 Data File : BP019269.D
 Acq On : 04 Jan 2024 18:14
 Operator : MA/JU
 Sample : 05951-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCF81

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-BP122723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP019269.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.899	129	138	143	rBV	65506	155899	0.49%	0.038%
2	5.287	368	374	386	rBV	33622	67933	0.22%	0.016%
3	5.516	406	413	421	rBV	126204	185259	0.59%	0.045%
4	6.093	505	511	516	rVB	60576	87433	0.28%	0.021%
5	6.699	610	614	623	rVB	46443	75041	0.24%	0.018%
6	6.840	632	638	644	rBV	51476	82525	0.26%	0.020%
7	6.993	659	664	675	rVB	142283	257121	0.82%	0.062%
8	7.134	681	688	700	rBV	131369	231175	0.73%	0.056%
9	7.334	715	722	731	rBV	161347	278408	0.88%	0.067%
10	7.434	731	739	745	rVB	133659	219201	0.70%	0.053%
11	7.552	755	759	768	rVB3	36584	69857	0.22%	0.017%
12	7.793	791	800	806	rBV	742549	1177226	3.74%	0.284%
13	7.869	806	813	826	rVV	2124423	3492049	11.08%	0.843%
14	8.410	899	905	920	rVB4	36837	101666	0.32%	0.025%
15	8.534	920	926	932	rBV3	68905	137227	0.44%	0.033%
16	8.963	993	999	1004	rBV	149431	257609	0.82%	0.062%
17	9.016	1004	1008	1017	rVB	58477	106962	0.34%	0.026%
18	9.122	1017	1026	1035	rBV	201202	447694	1.42%	0.108%
19	9.422	1072	1077	1082	rVB3	38300	72594	0.23%	0.018%
20	9.487	1082	1088	1092	rBV3	55327	122926	0.39%	0.030%
21	9.681	1115	1121	1126	rBV3	134616	229519	0.73%	0.055%
22	9.898	1150	1158	1162	rBV3	54714	132635	0.42%	0.032%
23	9.993	1171	1174	1180	rVB3	57769	105925	0.34%	0.026%
24	10.110	1189	1194	1200	rBV	174260	331628	1.05%	0.080%
25	10.169	1202	1204	1209	rVB3	67897	94025	0.30%	0.023%
26	10.234	1209	1215	1220	rBV	169868	271499	0.86%	0.066%
27	10.298	1220	1226	1233	rVB3	68308	167112	0.53%	0.040%
28	10.369	1233	1238	1241	rBV4	57832	95696	0.30%	0.023%
29	10.481	1249	1257	1259	rBV4	67058	142857	0.45%	0.035%
30	10.563	1265	1271	1274	rBV	1362917	2343764	7.44%	0.566%
31	10.593	1274	1276	1281	rVB	932499	1196506	3.80%	0.289%
32	10.704	1291	1295	1299	rBV2	393930	597880	1.90%	0.144%
33	10.746	1299	1302	1306	rVB	235036	302932	0.96%	0.073%
34	10.816	1309	1314	1320	rBV2	495758	825118	2.62%	0.199%
35	10.875	1320	1324	1327	rBV2	213496	330922	1.05%	0.080%
36	10.916	1327	1331	1336	rVB	628864	905845	2.87%	0.219%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
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37	10.963	1336	1339	1342	rBV3	83331	120362	0.38%	0.029%
38	11.075	1353	1358	1359	rBV3	135603	194126	0.62%	0.047%
39	11.104	1359	1363	1370	rVB	1002898	1498399	4.75%	0.362%
40	11.198	1375	1379	1384	rVB2	518569	740618	2.35%	0.179%
41	11.251	1384	1388	1390	rBV	152812	233397	0.74%	0.056%
42	11.428	1412	1418	1422	rBV5	348017	746935	2.37%	0.180%
43	11.504	1428	1431	1438	rVB3	366097	661229	2.10%	0.160%
44	11.610	1444	1449	1454	rBV	819542	1614029	5.12%	0.390%
45	11.698	1461	1464	1470	rVB3	263617	441406	1.40%	0.107%
46	11.857	1484	1491	1497	rVB4	581617	1382045	4.39%	0.334%
47	11.922	1499	1502	1505	rBV2	166483	228909	0.73%	0.055%
48	12.016	1508	1518	1529	rBV5	1948739	6824522	21.65%	1.648%
49	12.145	1538	1540	1545	rVB4	231946	285986	0.91%	0.069%
50	12.216	1545	1552	1554	rVV6	461291	1000124	3.17%	0.242%
51	12.269	1558	1561	1565	rVB3	268291	352271	1.12%	0.085%
52	12.381	1576	1580	1583	rBV2	371817	505424	1.60%	0.122%
53	12.481	1589	1597	1607	rBV6	466181	1518081	4.82%	0.367%
54	12.675	1620	1630	1640	rBV4	1890062	4679575	14.85%	1.130%
55	12.757	1640	1644	1647	rBV3	562158	726039	2.30%	0.175%
56	12.828	1652	1656	1659	rBV2	716846	1107482	3.51%	0.267%
57	12.916	1667	1671	1676	rBV2	983893	1604183	5.09%	0.387%
58	12.975	1676	1681	1686	rVV2	2512946	3801094	12.06%	0.918%
59	13.092	1697	1701	1704	rBV2	416149	702440	2.23%	0.170%
60	13.187	1714	1717	1720	rBV2	296134	336297	1.07%	0.081%
61	13.269	1722	1731	1737	rBV	3980925	7654913	24.29%	1.849%
62	13.328	1738	1741	1742	rBV3	271764	319272	1.01%	0.077%
63	13.357	1742	1746	1753	rBV3	642172	1889720	6.00%	0.456%
64	13.475	1762	1766	1767	rBV2	572437	691934	2.20%	0.167%
65	13.498	1767	1770	1774	rVB3	988676	1409385	4.47%	0.340%
66	13.610	1781	1789	1796	rBV2	2541864	7310511	23.20%	1.766%
67	13.769	1811	1816	1820	rBV	4554736	8452016	26.82%	2.041%
68	13.928	1835	1843	1846	rBV2	4501639	7987869	25.34%	1.929%
69	13.963	1846	1849	1853	rVV3	2362883	3491304	11.08%	0.843%
70	14.004	1853	1856	1859	rVV	1166887	1469980	4.66%	0.355%
71	14.039	1859	1862	1867	rVB2	2083051	3085842	9.79%	0.745%
72	14.092	1867	1871	1875	rBV4	875372	1533212	4.86%	0.370%
73	14.134	1875	1878	1881	rBV2	1264875	1790762	5.68%	0.432%
74	14.363	1909	1917	1922	rBV2	16084711	31516595	100.00%	7.612%
75	14.422	1923	1927	1928	rBV4	1366301	1817212	5.77%	0.439%
76	14.581	1951	1954	1957	rBV2	1571329	1838982	5.83%	0.444%

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 Title : SVOA CALIBRATION

77	14.681	1964	1971	1976	rVB3	4747363	13000975	41.25%	3.140%
78	14.751	1978	1983	1986	rBV2	1854582	3838469	12.18%	0.927%
79	14.857	1997	2001	2006	rVB2	8210370	12372762	39.26%	2.988%
80	14.934	2007	2014	2018	rBV2	6289157	11398609	36.17%	2.753%
81	14.975	2018	2021	2026	rVV5	4967983	7322832	23.23%	1.769%
82	15.045	2030	2033	2035	rVV	1955106	2674158	8.48%	0.646%
83	15.104	2035	2043	2046	rVV2	9970955	23822748	75.59%	5.753%
84	15.134	2046	2048	2053	rVB	4443515	4955693	15.72%	1.197%
85	15.286	2070	2074	2076	rBV2	5644222	7765695	24.64%	1.875%
86	15.328	2077	2081	2084	rVB	11039164	16157922	51.27%	3.902%
87	15.575	2116	2123	2128	rBV4	7675582	21242486	67.40%	5.130%
88	15.728	2145	2149	2153	rVB	9670242	15602936	49.51%	3.768%
89	15.881	2170	2175	2180	rBV4	5194720	10776538	34.19%	2.603%
90	15.939	2180	2185	2188	rBV3	5386785	9234509	29.30%	2.230%
91	16.204	2223	2230	2233	rBV2	11576259	30566693	96.99%	7.382%
92	16.916	2342	2351	2352	rBV3	11401310	29435572	93.40%	7.109%
93	19.027	2706	2710	2715	rBV2	6546612	13587922	43.11%	3.282%
94	20.574	2970	2973	2982	rVB2	8438342	13331944	42.30%	3.220%
95	21.027	3047	3050	3053	rVB2	5661420	6428840	20.40%	1.553%
96	21.227	3081	3084	3090	rVB	7373253	9425077	29.91%	2.276%
97	21.457	3120	3123	3131	rVB2	4496297	6953893	22.06%	1.679%
98	23.633	3488	3493	3498	rBV	1920821	3437340	10.91%	0.830%
99	25.145	3745	3750	3754	rBV2	1402165	3117211	9.89%	0.753%
100	25.404	3786	3794	3806	rVB	3025130	8347557	26.49%	2.016%

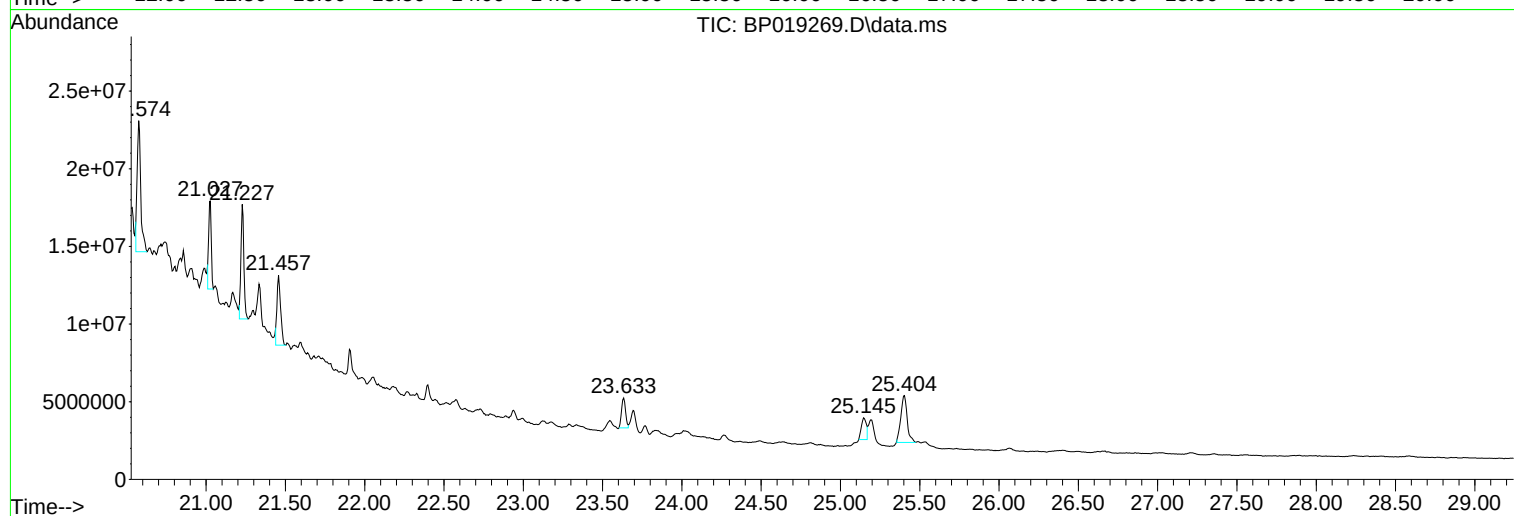
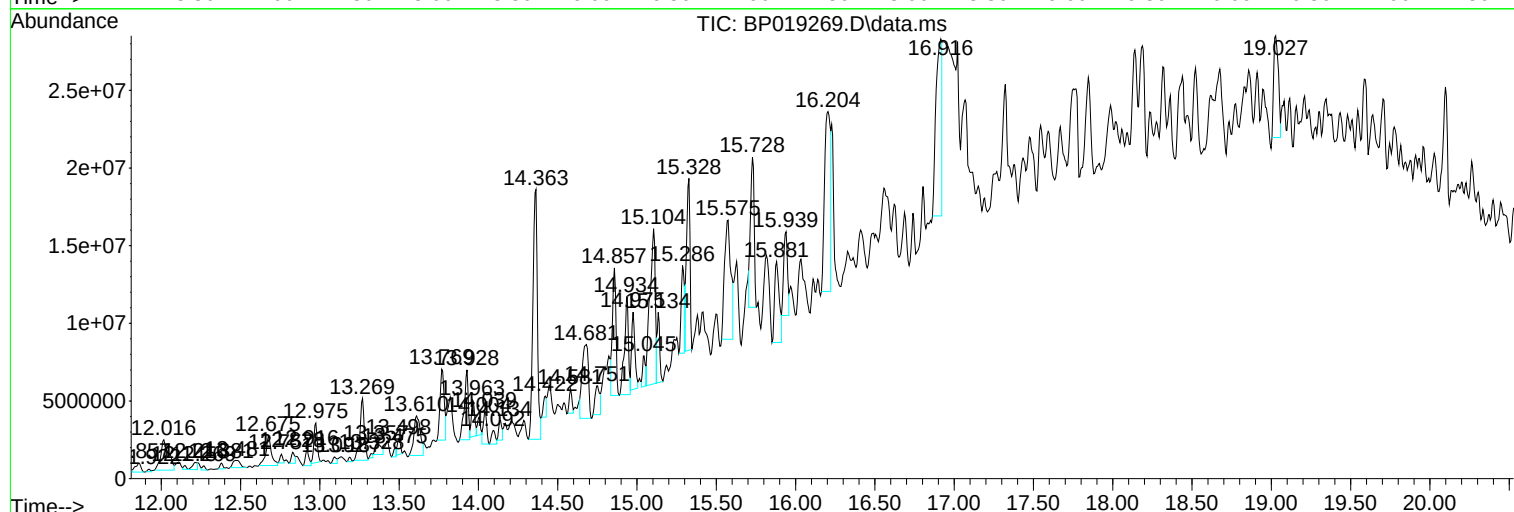
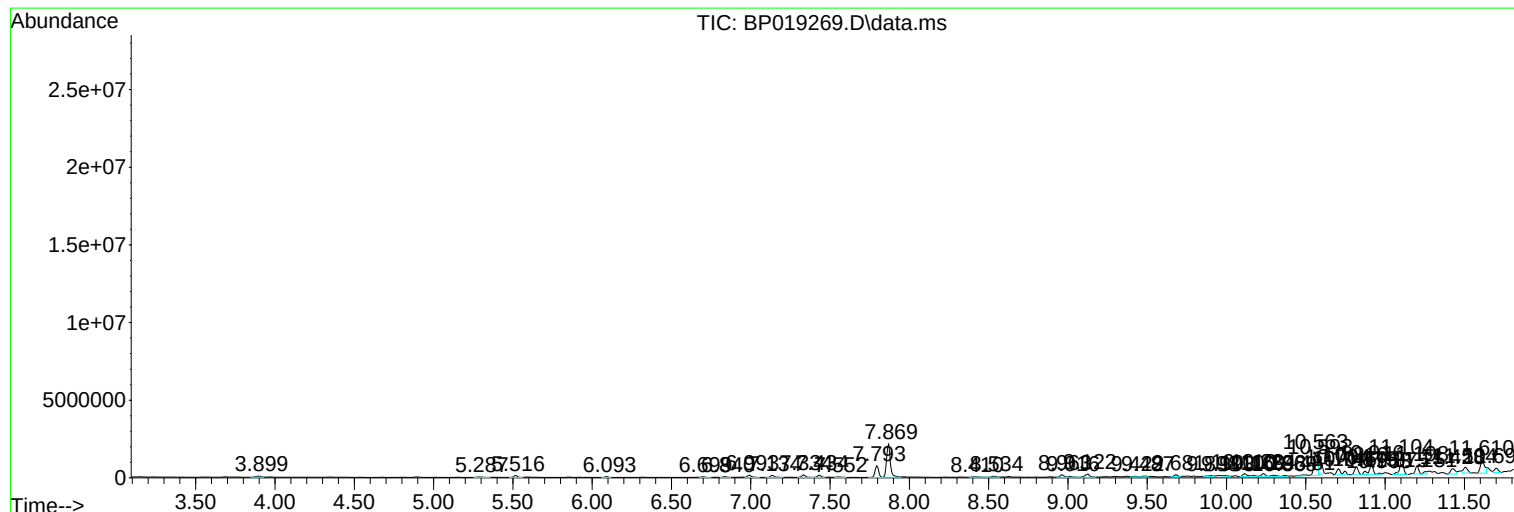
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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



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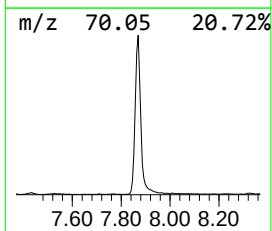
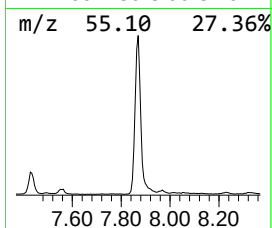
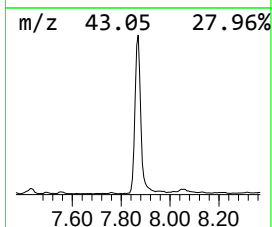
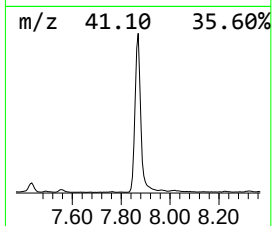
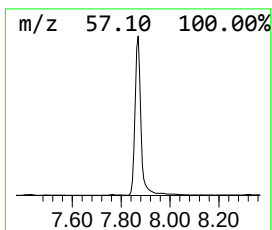
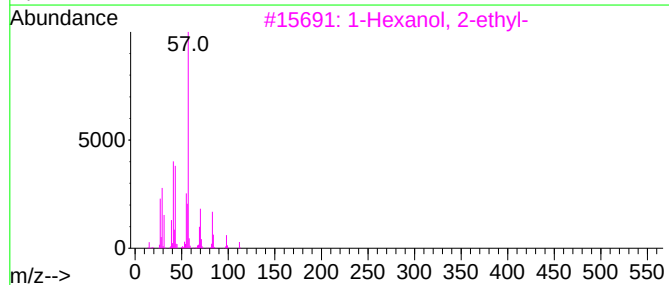
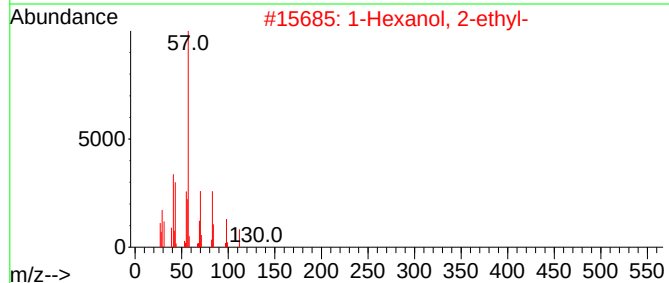
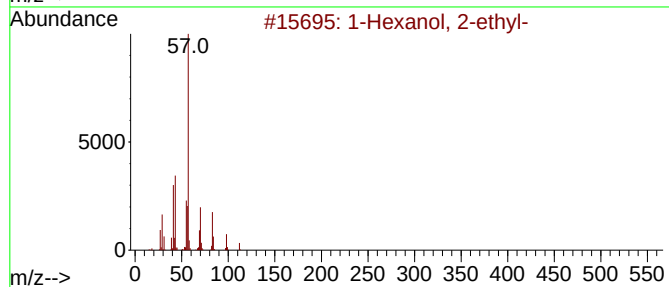
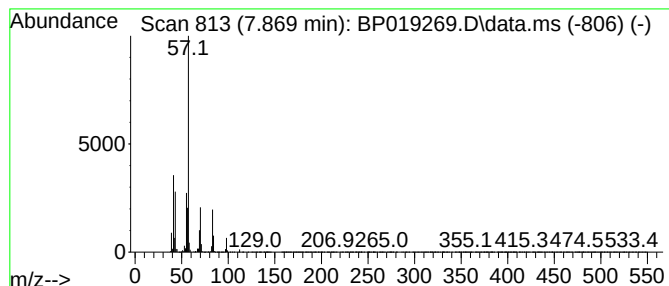
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Hexanol, 2-ethyl- Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.869	59.33 ng/ul	3492050	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	78
2	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	72
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	64
4	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	56
5	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	53



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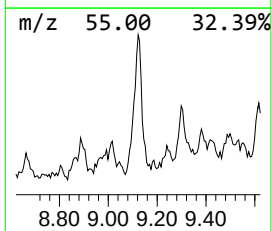
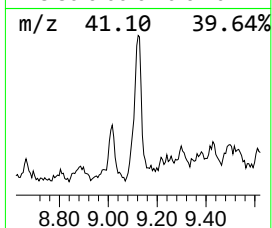
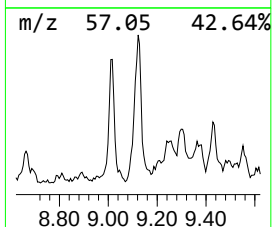
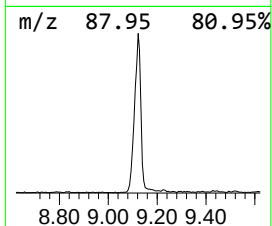
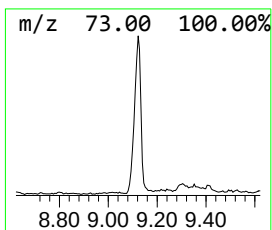
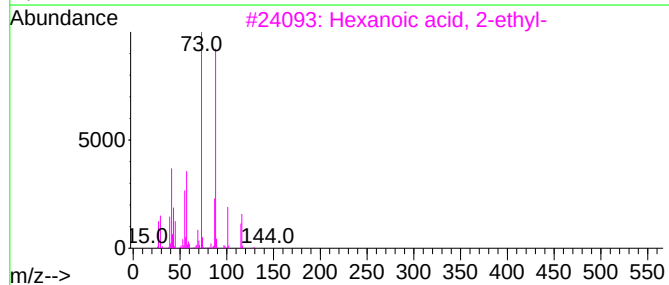
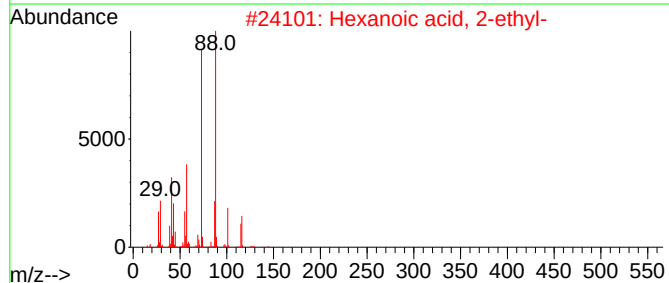
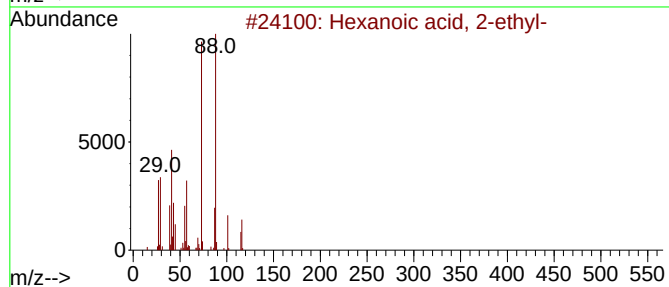
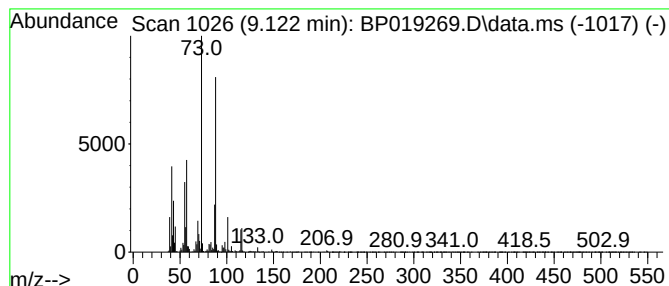
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Hexanoic acid, 2-ethyl- Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.122	7.61 ng/ul	447694	1,4-Dichlorobenzene-d4	7.799

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	86
2			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	83
3			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	80
4			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78
5			Hexanoic acid, 2-ethyl-	144	C8H16O2	000149-57-5	78



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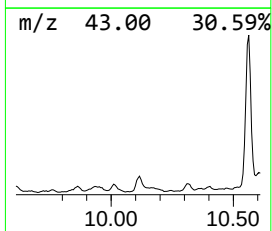
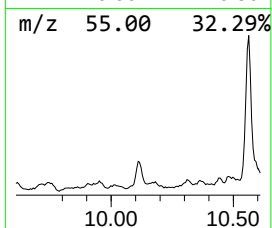
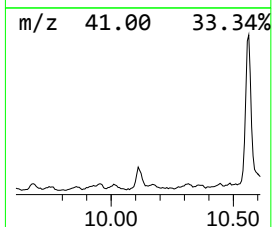
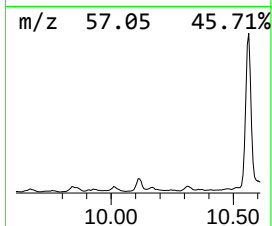
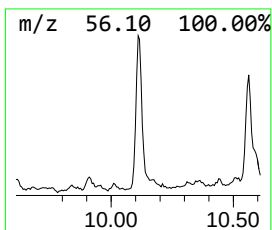
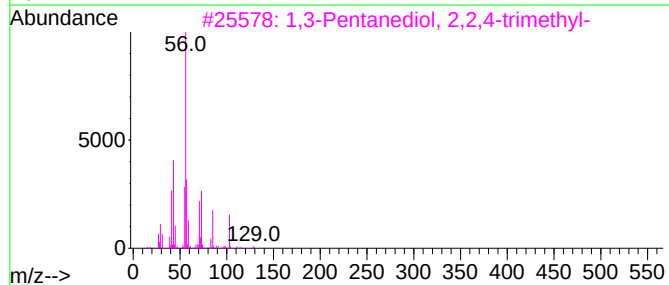
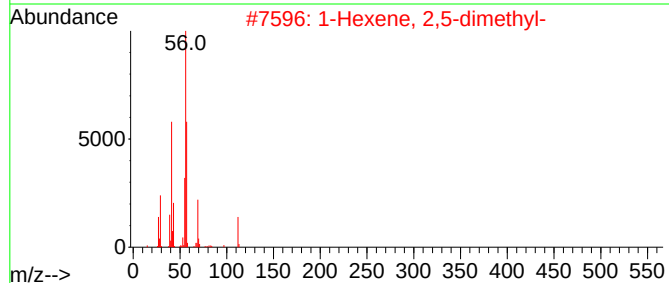
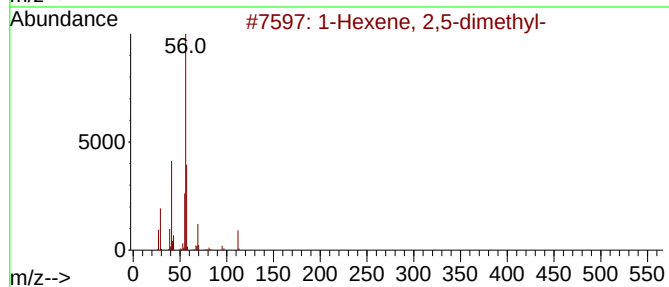
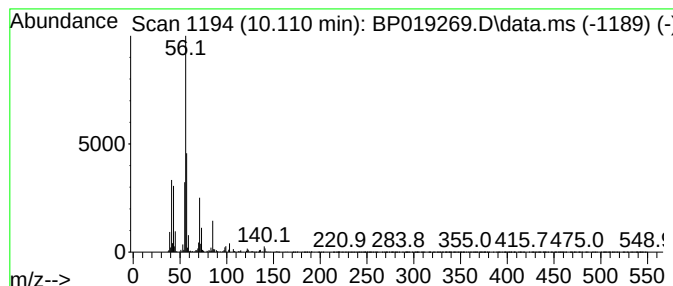
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 50

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.110	5.54 ng/ul	331628	Naphthalene-d8	10.593

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	43
2		1-Hexene, 2,5-dimethyl-	112	C8H16	006975-92-4	43
3		1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	38
4		1-Hexene, 2-methyl-	98	C7H14	006094-02-6	38
5		1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	37



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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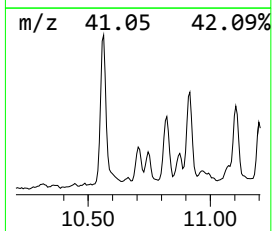
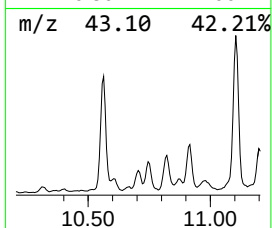
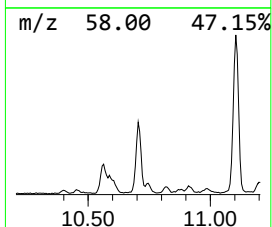
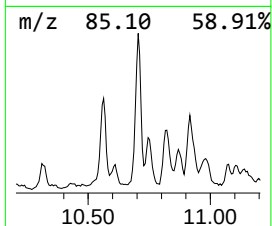
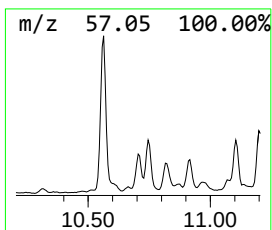
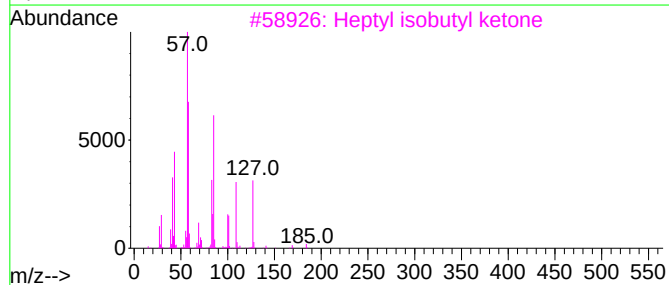
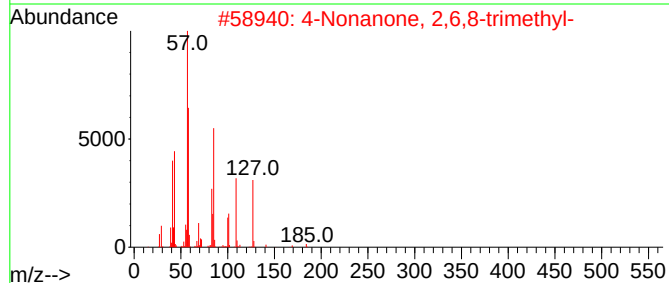
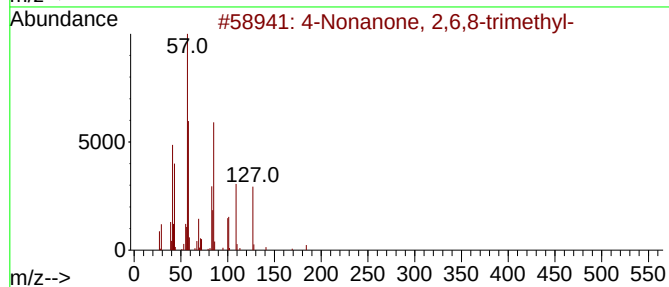
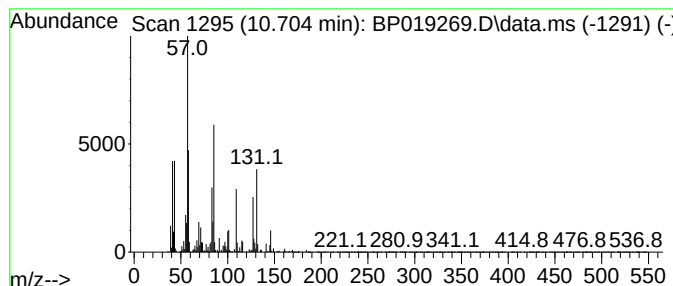
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 4-Nonanone, 2,6,8-trimethyl- Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.704	9.99 ng/ul	597880	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	95
2			4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	90
3			Heptyl isobutyl ketone	184	C12H24O	019594-40-2	46
4			4-Octanone, 2-methyl-	142	C9H18O	007492-38-8	27
5			3,3-Dimethylbutan-2-yl isobutyl ...	202	C11H22O3	1000372-76-5	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
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Sample : 05951-17
Misc :
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Instrument :
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ClientSampleId :
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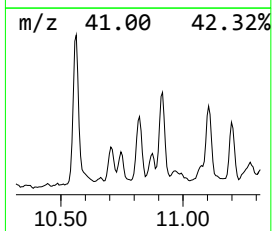
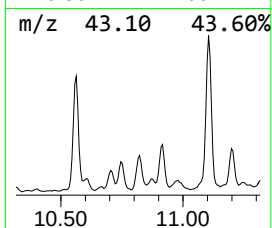
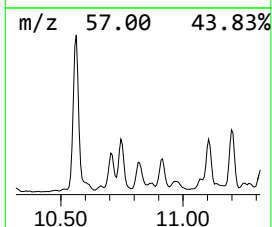
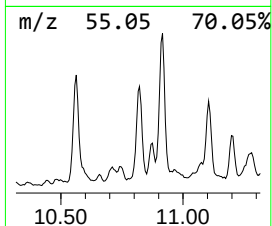
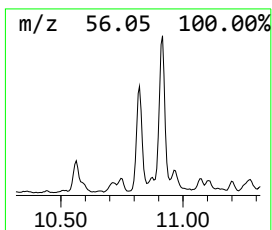
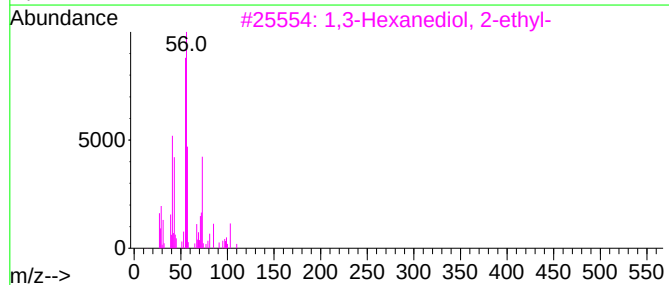
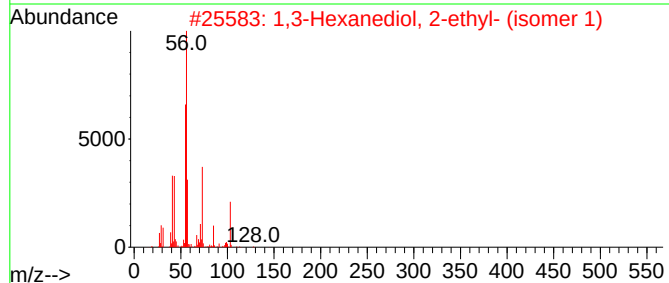
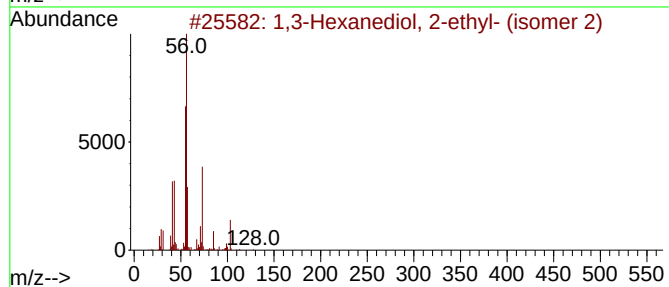
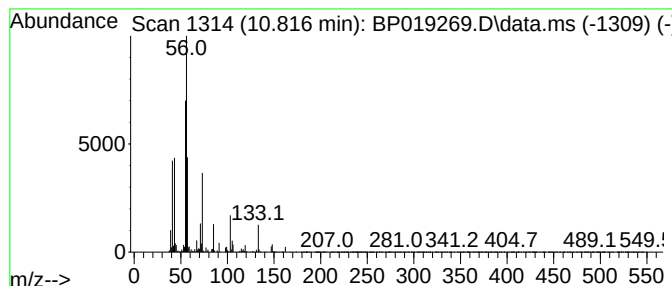
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 1,3-Hexanediol, 2-ethyl- (i... Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	13.79 ng/ul	825118	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Hexanediol, 2-ethyl- (isomer 2)	146	C8H18O2	1000484-18-1	72		
2	1,3-Hexanediol, 2-ethyl- (isomer 1)	146	C8H18O2	1000484-18-2	64		
3	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	56		
4	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	47		
5	1,3-Pentanediol, 2,2,4-trimethyl-	146	C8H18O2	000144-19-4	45		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

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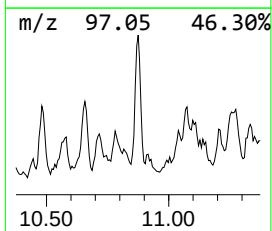
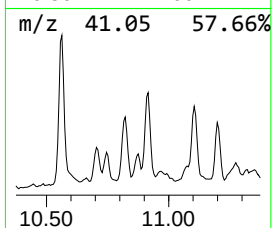
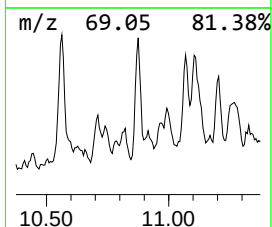
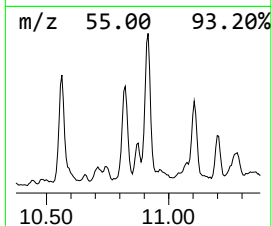
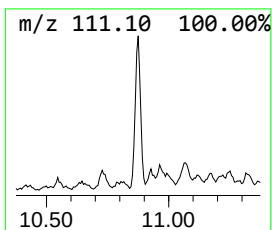
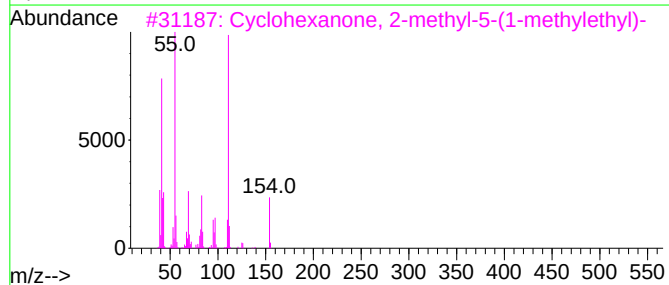
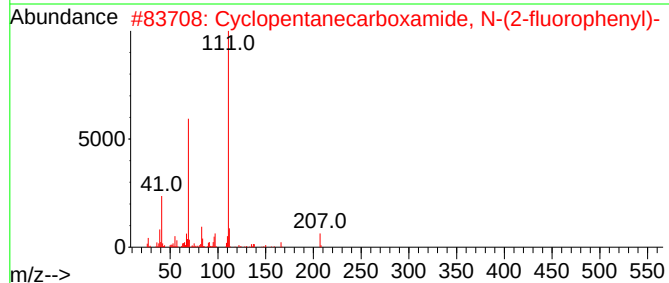
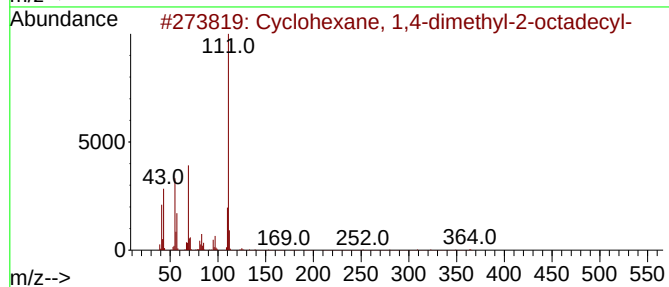
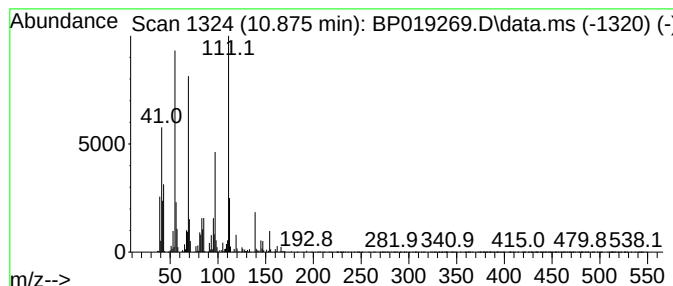
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Cyclic10.875 Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.875	5.53 ng/ul	330922	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, 1,4-dimethyl-2-octa...	364	C26H52	055282-02-5	58
2			Cyclopentanecarboxamide, N-(2-fl...	207	C12H14FNO	1000308-60-7	49
3			Cyclohexanone, 2-methyl-5-(1-met...	154	C10H18O	059471-80-6	41
4			3-Heptene, 2-methyl-, (E)-	112	C8H16	000692-96-6	38
5			3-Heptene, 4-methyl-	112	C8H16	004485-16-9	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
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Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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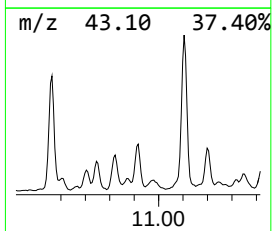
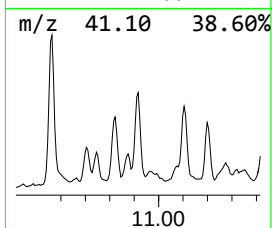
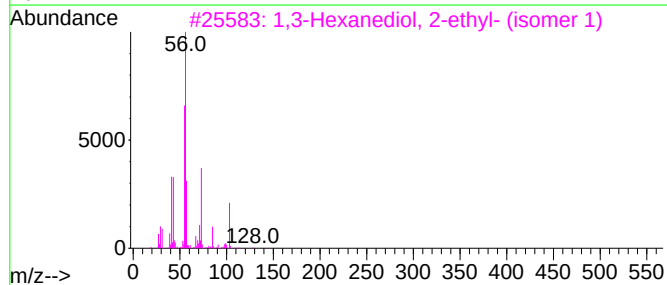
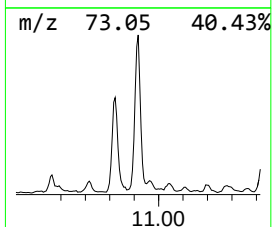
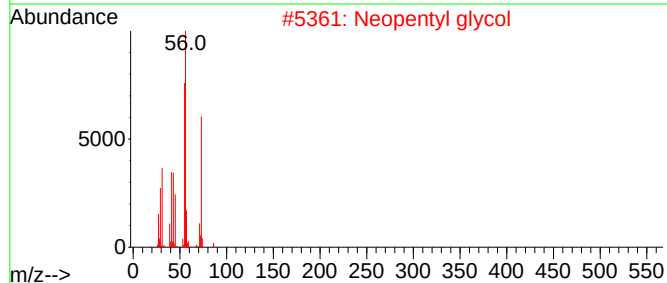
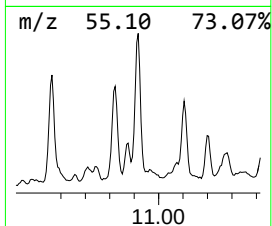
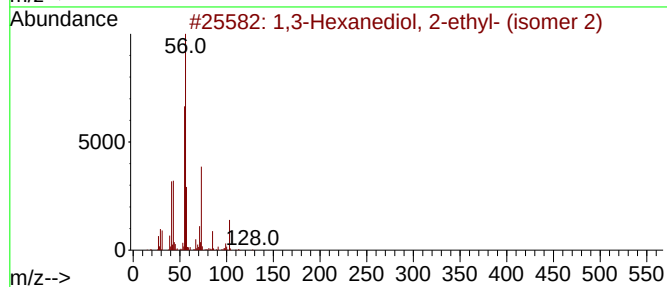
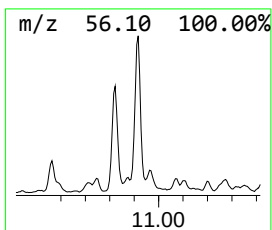
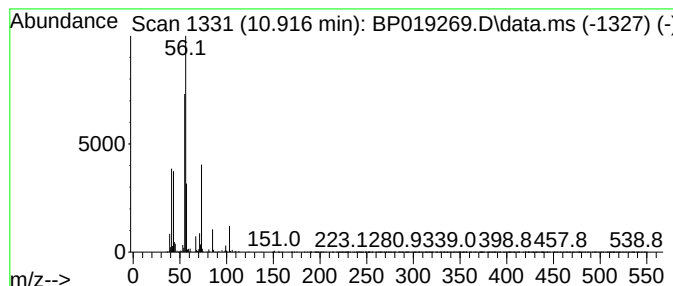
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Neopentyl glycol Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.916	15.14 ng/ul	905845	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,3-Hexanediol, 2-ethyl- (isomer 2)	146	C8H18O2	1000484-18-1	90		
2	Neopentyl glycol	104	C5H12O2	000126-30-7	64		
3	1,3-Hexanediol, 2-ethyl- (isomer 1)	146	C8H18O2	1000484-18-2	64		
4	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	64		
5	1,3-Hexanediol, 2-ethyl-	146	C8H18O2	000094-96-2	53		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
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Sample : 05951-17
Misc :
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Instrument :
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ClientSampleId :
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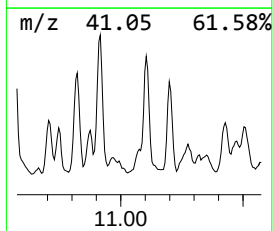
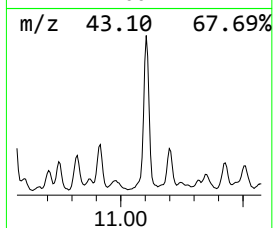
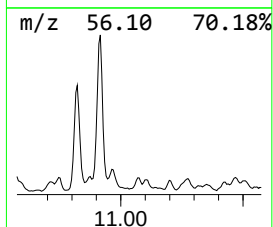
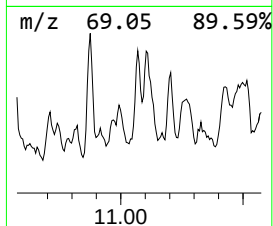
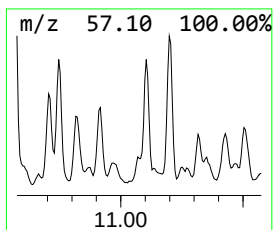
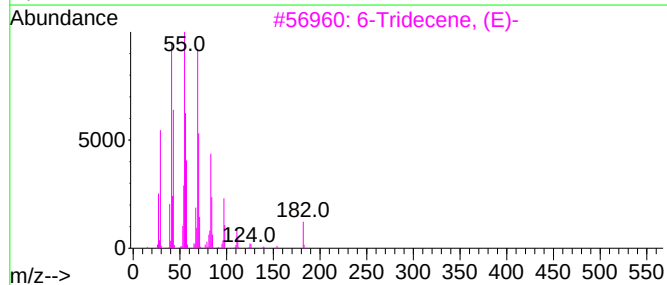
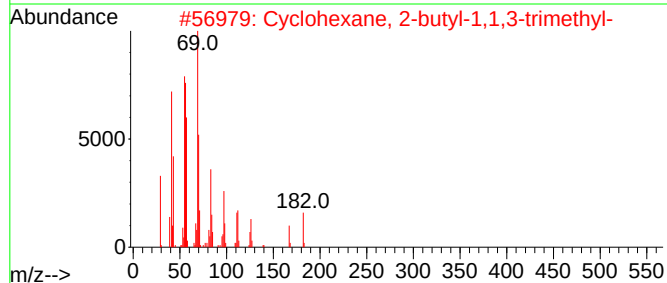
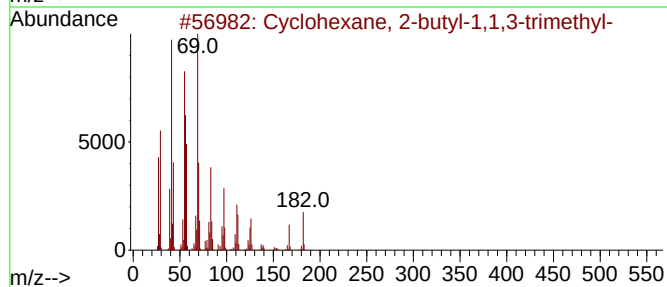
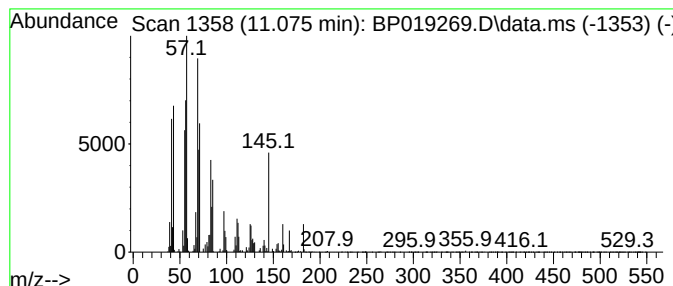
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 (DEL) Alkane: Cyclic11.075 Concentration Rank 59

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.075	3.24 ng/ul	194126	Naphthalene-d8	10.593

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	87
2		Cyclohexane, 2-butyl-1,1,3-trime...	182	C13H26	054676-39-0	78
3		6-Tridecene, (E)-	182	C13H26	006434-76-0	70
4		3-Tridecene, (E)-	182	C13H26	041446-57-5	70
5		3-Tridecene, (Z)-	182	C13H26	041446-53-1	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

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ClientSampleId :
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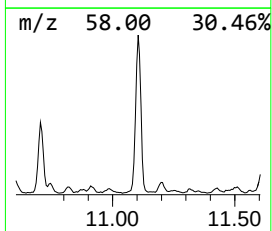
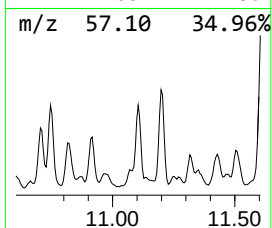
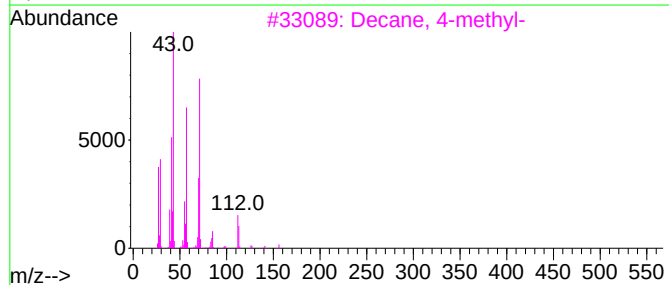
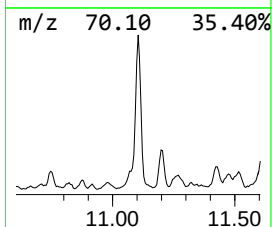
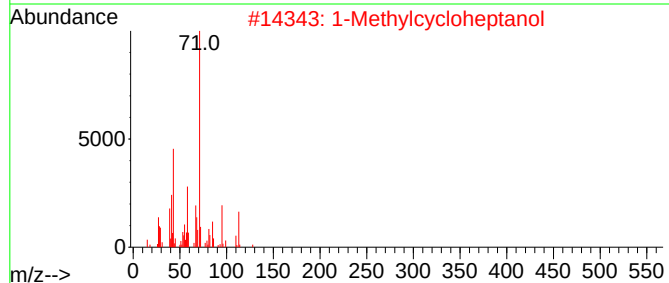
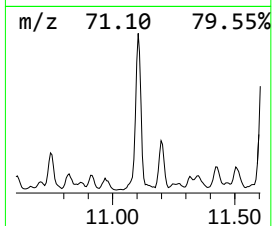
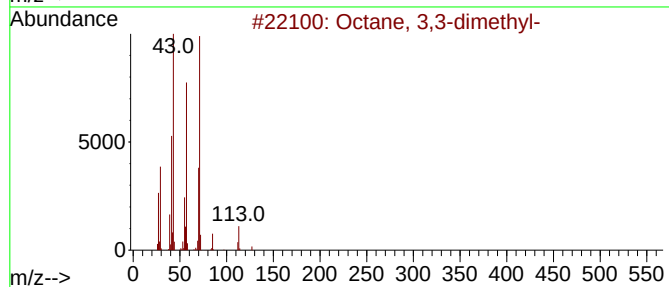
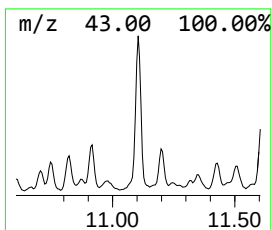
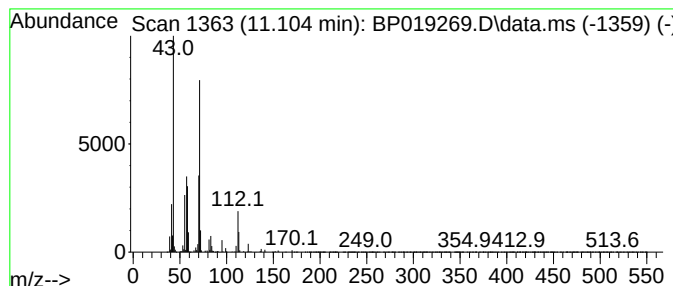
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.104	25.05 ng/ul	1498400	Naphthalene-d8	10.593

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	59	
2	1-Methylcycloheptanol	128	C8H16O	003761-94-2	50	
3	Decane, 4-methyl-	156	C11H24	002847-72-5	45	
4	1-Butanol, 2,2-dimethyl-	102	C6H14O	001185-33-7	42	
5	3-Acetyl-2-octanone	170	C10H18O2	027970-50-9	40	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
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Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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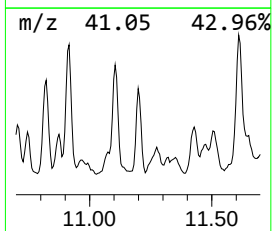
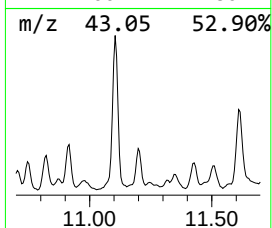
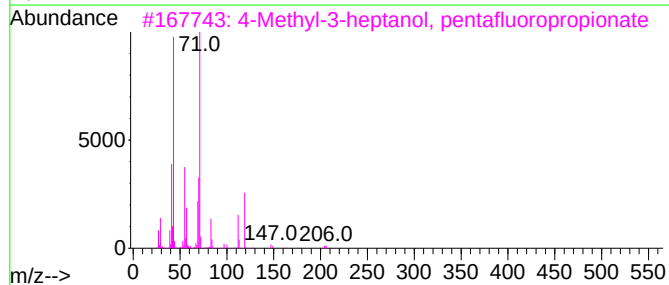
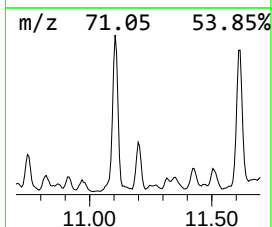
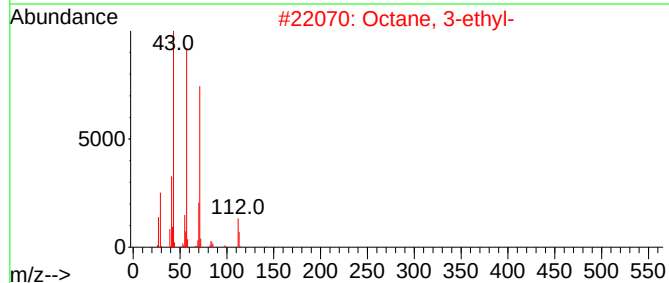
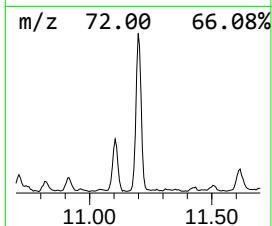
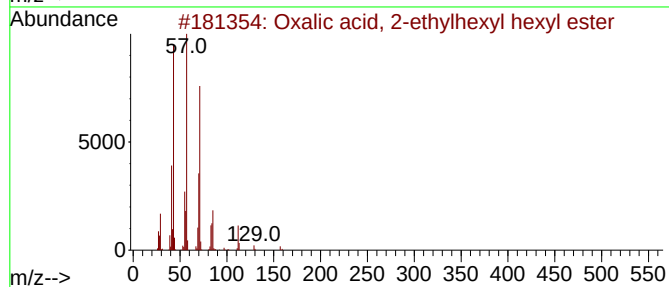
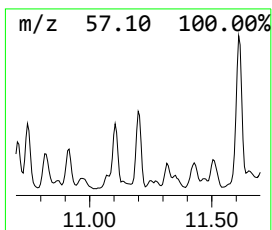
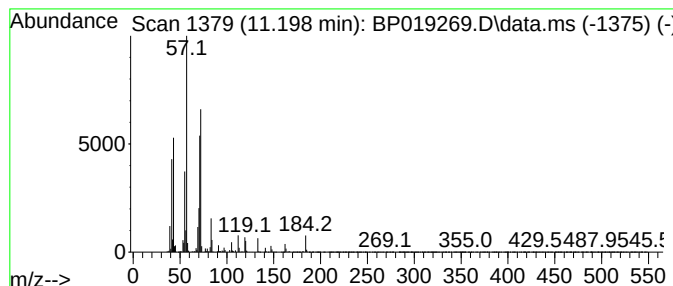
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 unknown-01 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.198	12.38 ng/ul	740618	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, 2-ethylhexyl hexyl ...	286	C16H30O4	1000309-38-9	35
2			Octane, 3-ethyl-	142	C10H22	005881-17-4	35
3			4-Methyl-3-heptanol, pentafluoro...	276	C11H17F5O2	1000365-51-7	32
4			Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	30
5			Propanal, 2-propenylhydrazone	112	C6H12N2	019031-78-8	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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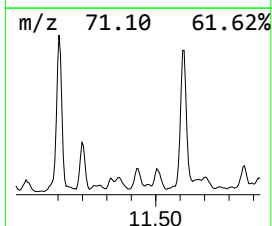
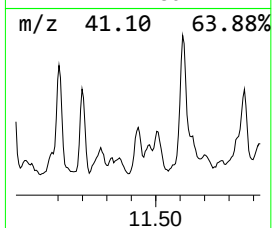
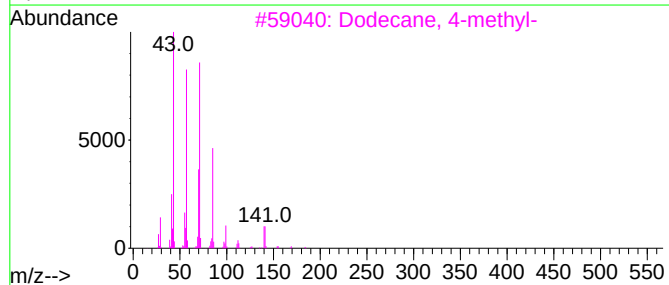
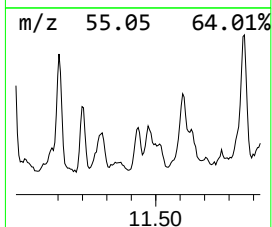
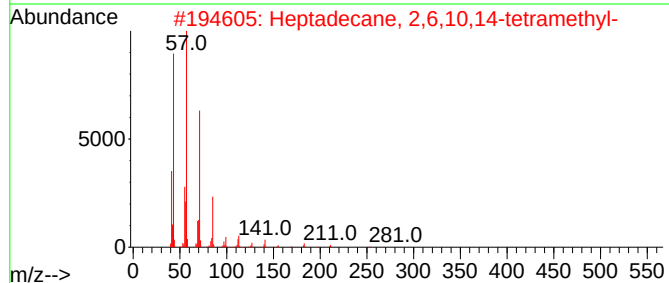
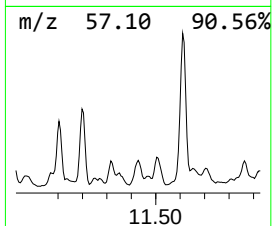
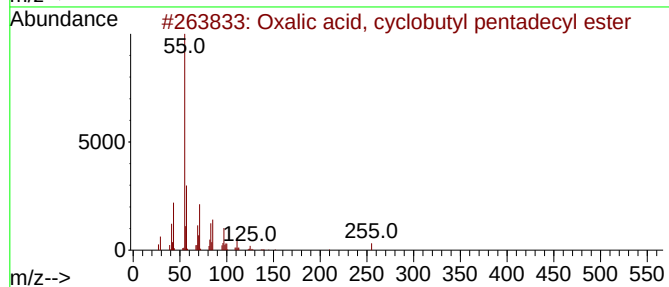
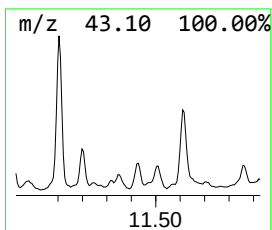
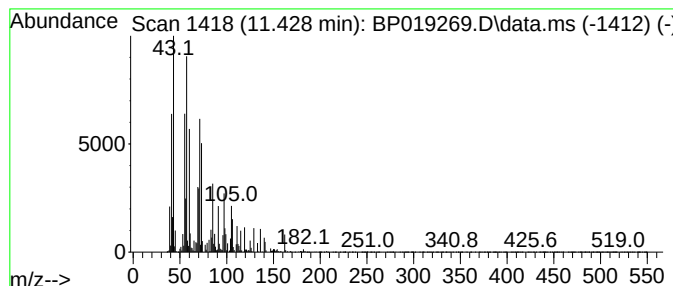
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 unknown-02 Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.428	12.49 ng/ul	746935	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, cyclobutyl pentadec...	354	C21H38O4	1010309-70-5	38
2			Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	35
3			Dodecane, 4-methyl-	184	C13H28	006117-97-1	35
4			Nonyl tetradecyl ether	340	C23H48O	1000406-37-6	30
5			Oxalic acid, allyl octadecyl ester	382	C23H42O4	1000309-24-5	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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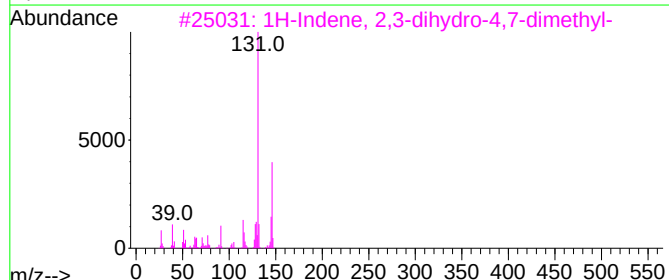
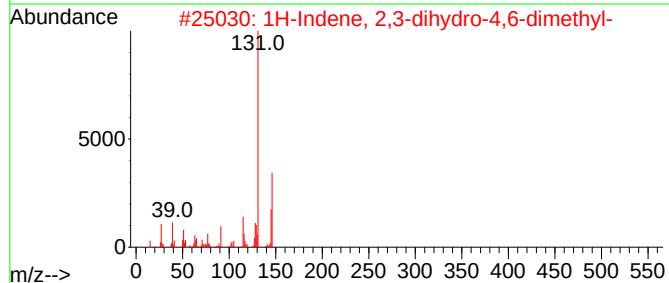
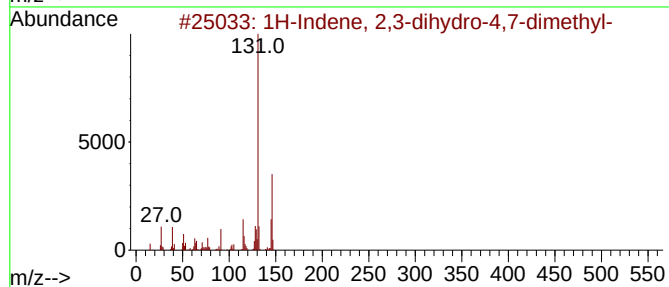
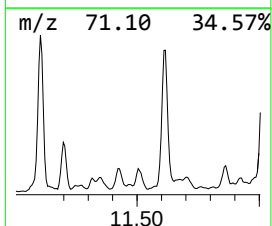
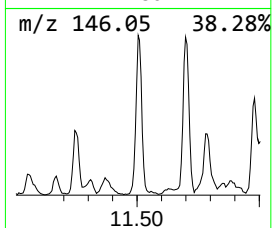
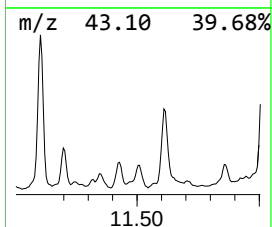
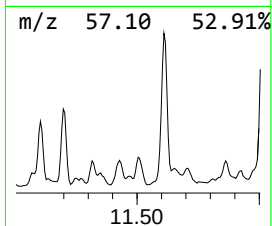
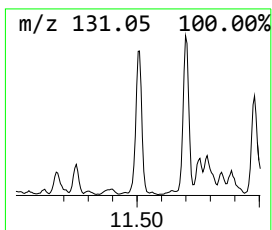
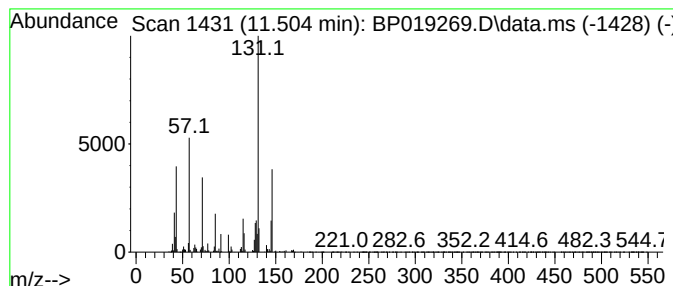
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 1H-Indene, 2,3-dihydro-4,7-... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.504	11.05 ng/ul	661229	Naphthalene-d8	10.593

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-4,6-dimet...	146	C11H14	001685-82-1	93		
3	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	91		
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	87		
5	1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	87		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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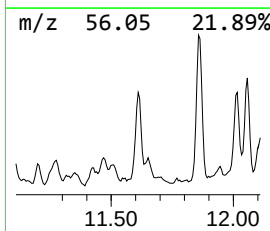
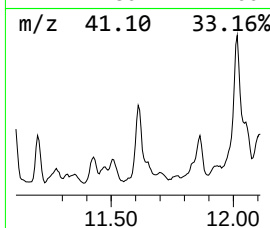
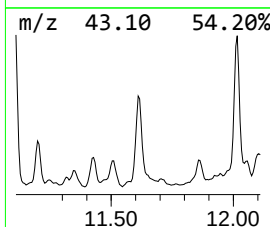
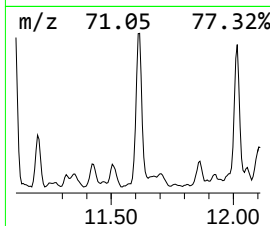
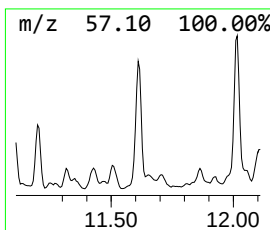
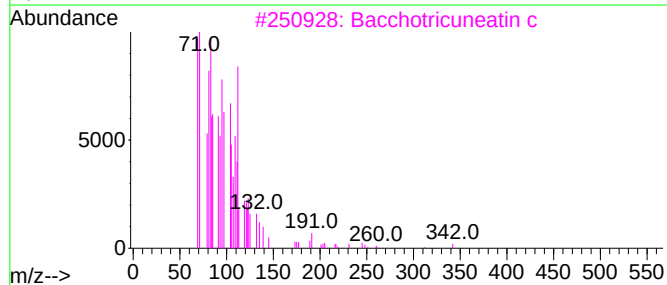
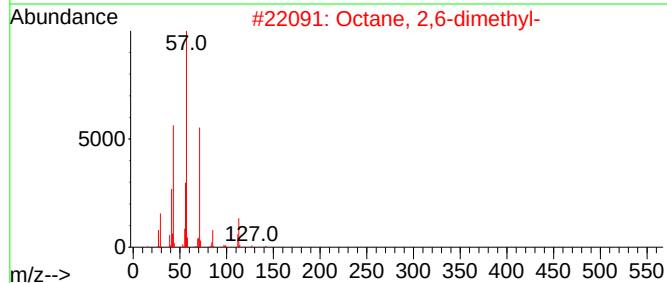
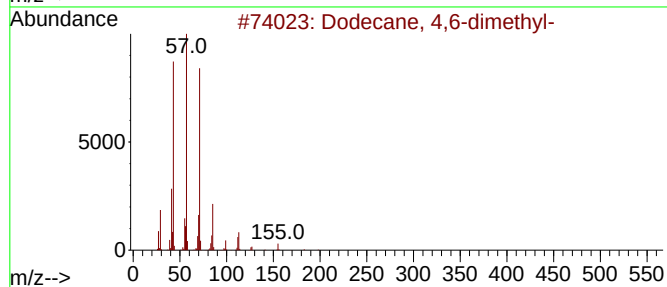
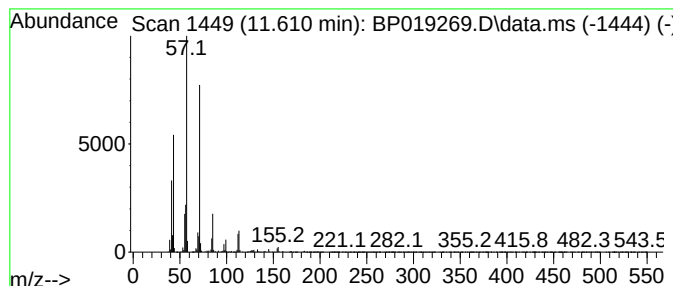
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 (DEL) Alkane: Straight-Chai... Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.610	26.98 ng/ul	1614030	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	87
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	86
3			Bacchotricuneatin c	342	C20H22O5	066563-30-2	72
4			Dodecane, 4,6-dimethyl-	198	C14H30	061141-72-8	72
5			Sulfurous acid, 2-ethylhexyl pen...	264	C13H28O3S	1000309-18-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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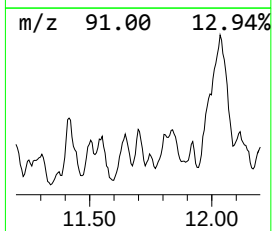
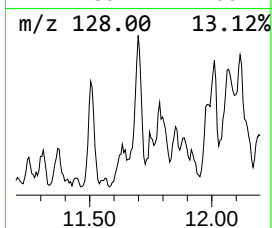
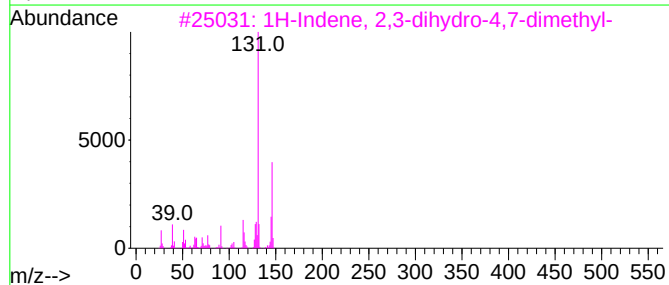
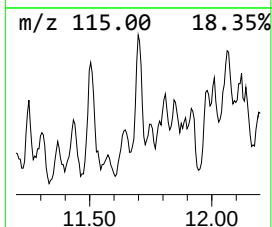
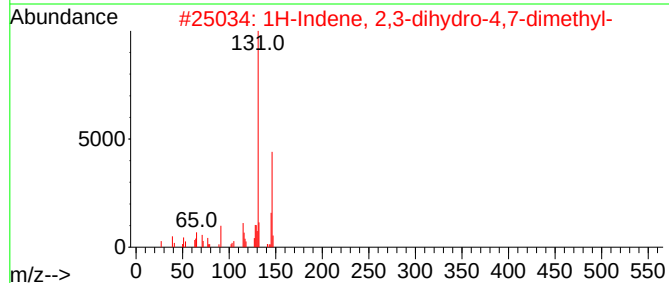
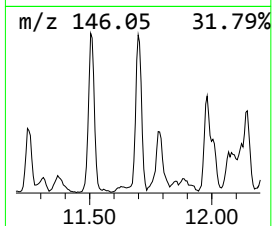
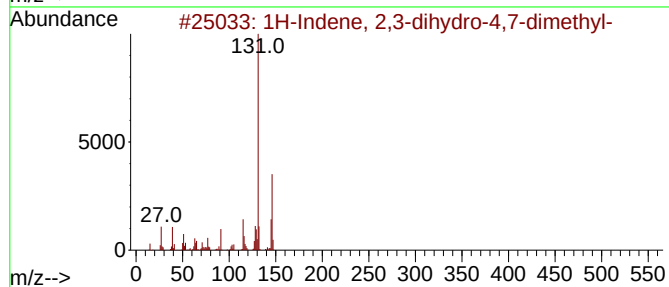
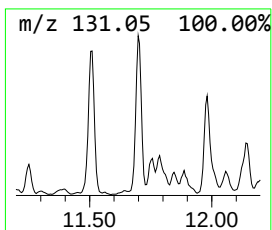
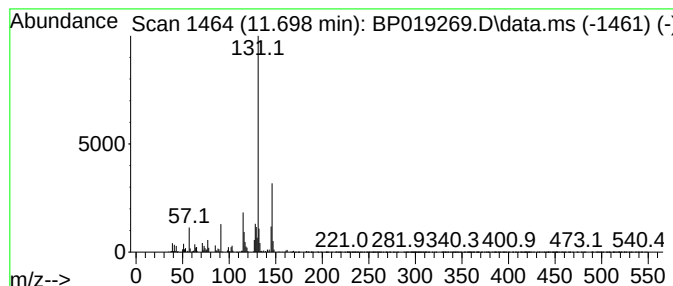
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Quant Title : SVOA CALIBRATION

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

Peak Number 24 Benzene, 2-ethenyl-1,3,5-tr... Concentration Rank 47

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.698	7.38 ng/ul	441406	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	97
2			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
3			1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	94
4			Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	92
5			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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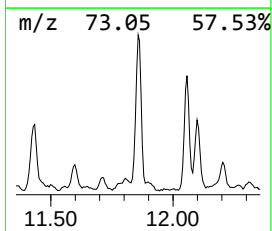
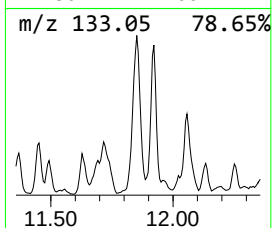
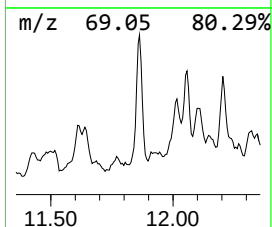
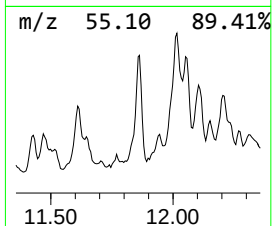
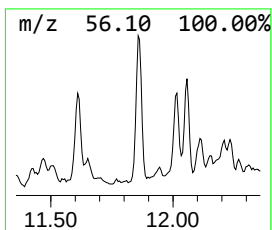
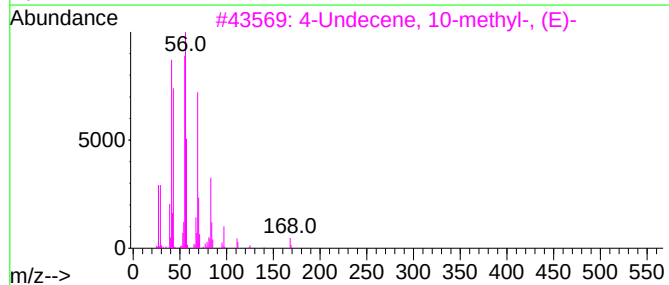
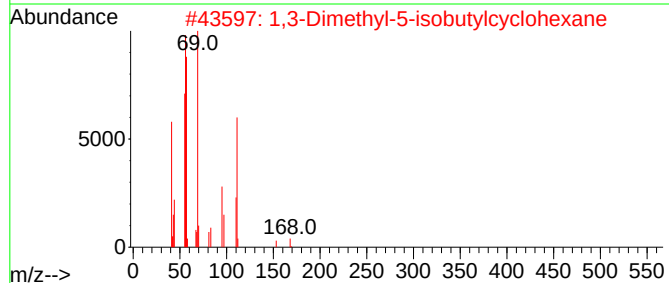
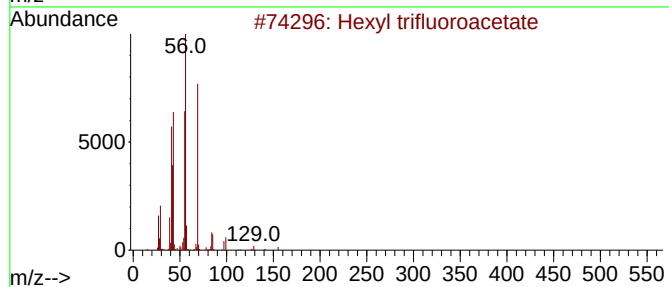
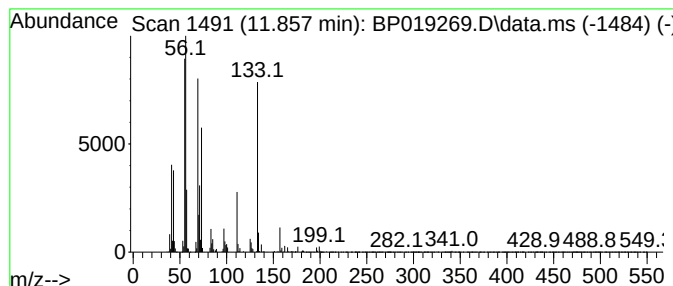
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Quant Title : SVOA CALIBRATION

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

Peak Number 25 unknown-03 Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.857	23.10 ng/ul	1382050	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexyl trifluoroacetate	198	C8H13F3O2	000400-61-3	38
2			1,3-Dimethyl-5-isobutylcyclohexane	168	C12H24	013131-76-5	37
3			4-Undecene, 10-methyl-, (E)-	168	C12H24	074630-60-7	37
4			2-Oxepanone, 4-methyl-	128	C7H12O2	002549-60-2	32
5			Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	32



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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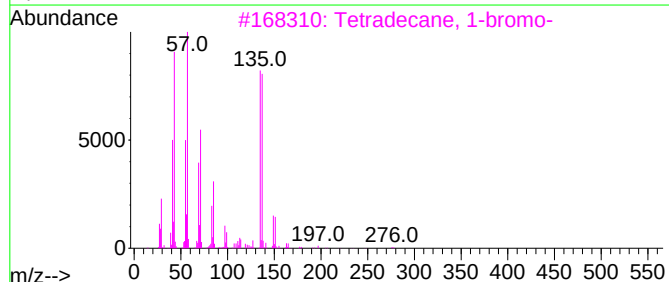
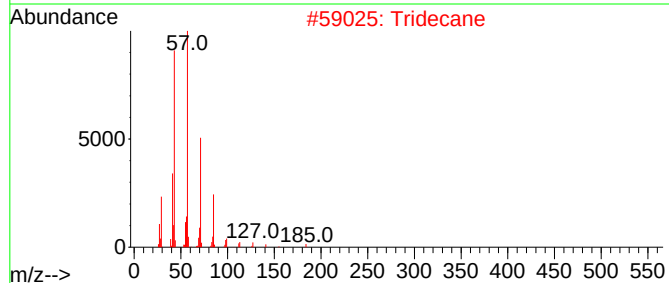
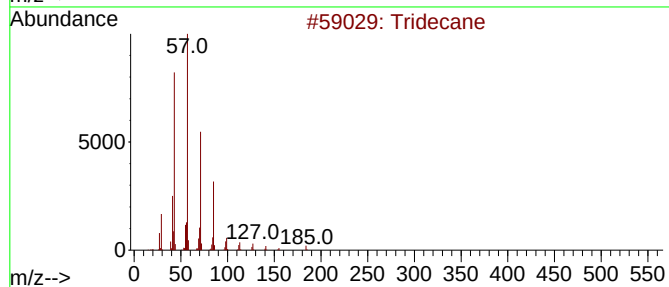
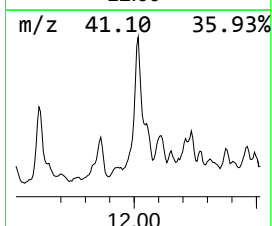
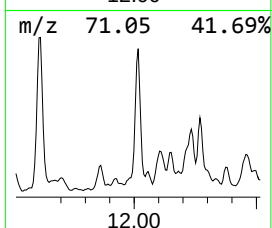
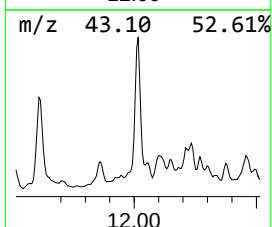
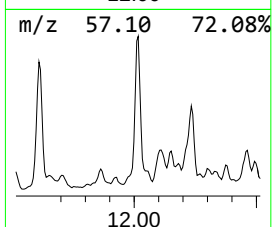
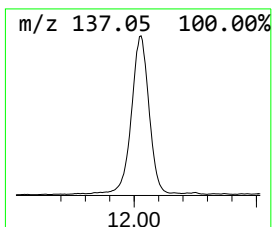
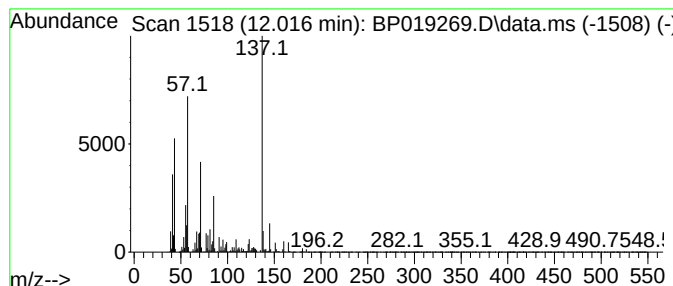
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Quant Title : SVOA CALIBRATION

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

Peak Number 27 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.016	114.07 ng/ul	6824520	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	55
2			Tridecane	184	C13H28	000629-50-5	38
3			Tetradecane, 1-bromo-	276	C14H29Br	000112-71-0	38
4			2,4-Dihydroxyphenylbenzyl ketone	228	C14H12O3	003669-41-8	38
5			Dodecane	170	C12H26	000112-40-3	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

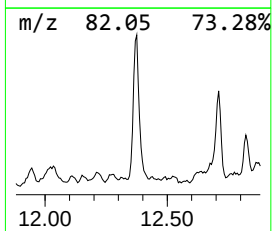
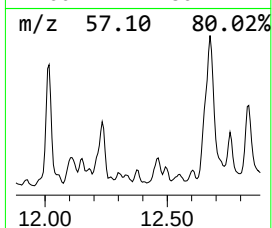
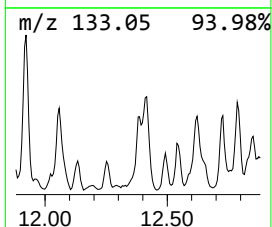
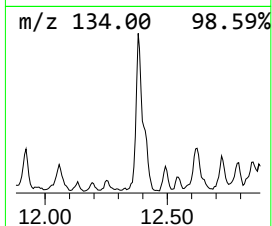
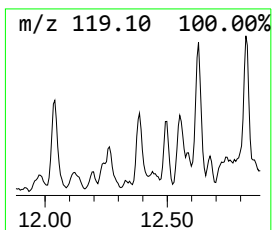
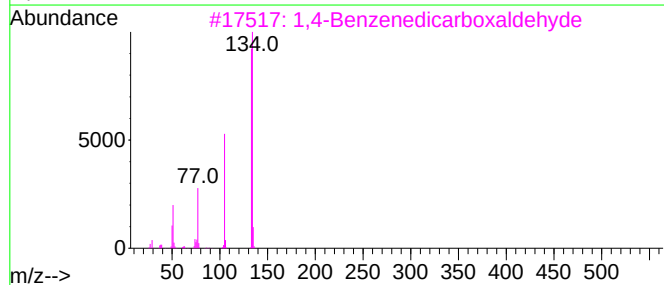
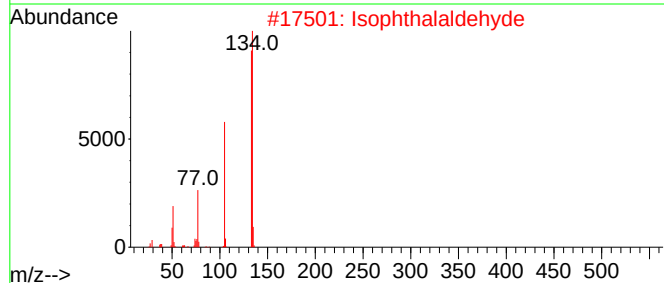
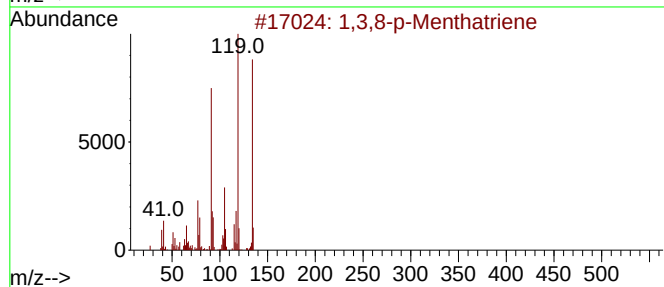
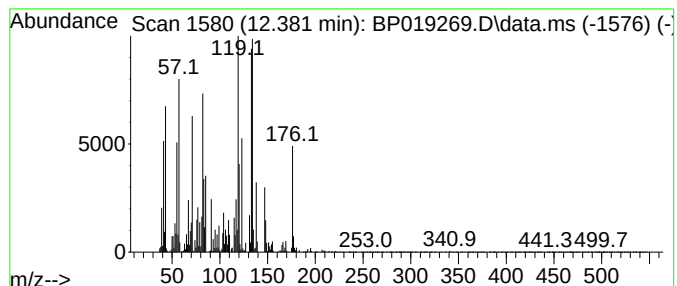
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TIC Integration Parameters: LSCINT.P

Peak Number 29 unknown-04

Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.381	8.45 ng/ul	505424	Naphthalene-d8	10.593

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3,8-p-Menthatriene	134	C10H14	018368-95-1	38
2			Isophthalaldehyde	134	C8H6O2	000626-19-7	35
3			1,4-Benzenedicarboxaldehyde	134	C8H6O2	000623-27-8	35
4			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	25
5			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
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Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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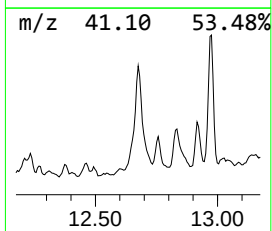
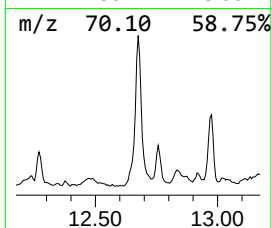
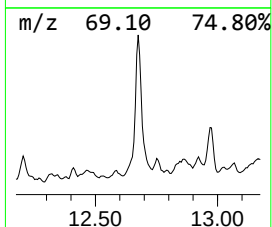
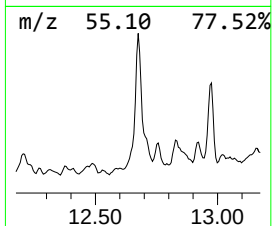
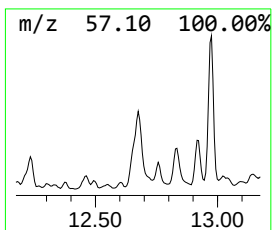
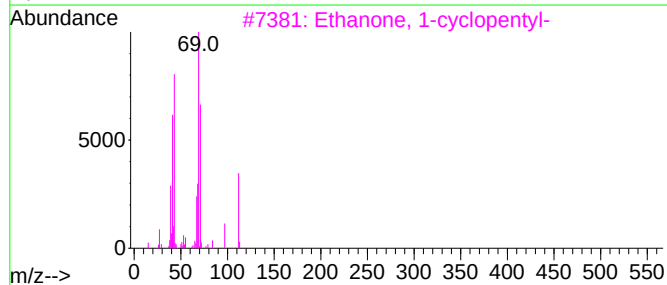
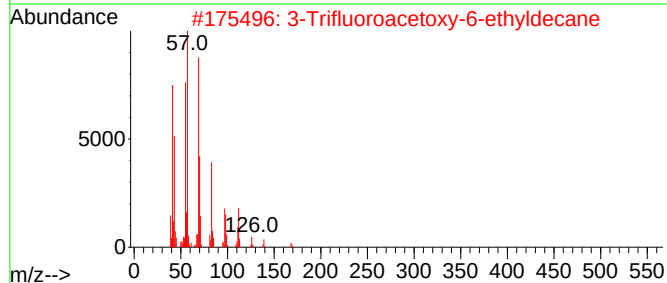
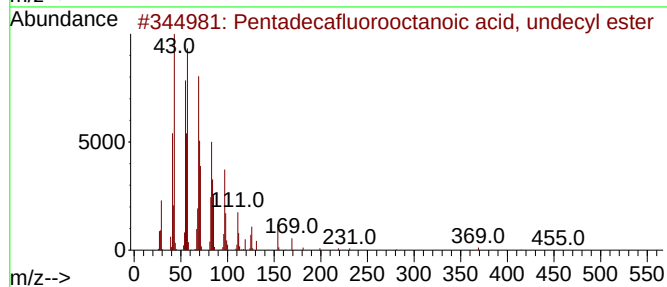
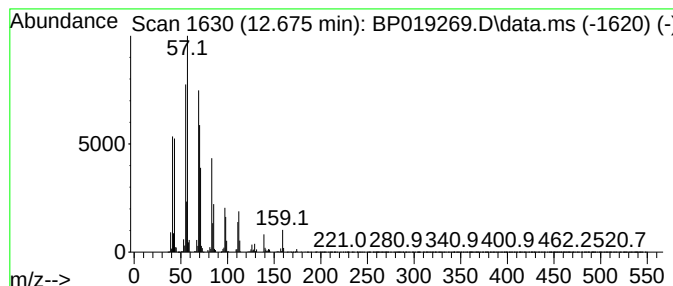
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Quant Title : SVOA CALIBRATION

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

Peak Number 30 Pentadecafluorooctanoic aci... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.675	51.50 ng/ul	4679580	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentadecafluorooctanoic acid, un...	568	C19H23F15O2	1010406-04-1	72
2			3-Trifluoroacetoxy-6-ethyldecane	282	C14H25F3O2	116436-59-0	49
3			Ethanone, 1-cyclopentyl-	112	C7H12O	006004-60-0	38
4			3-Hexen-2-one, 3-methyl-	112	C7H12O	001187-80-0	30
5			3-Heptene, 4-methyl-	112	C8H16	004485-16-9	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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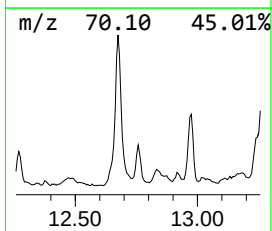
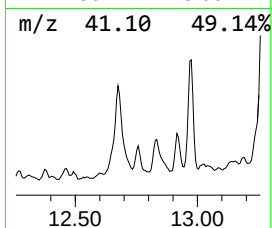
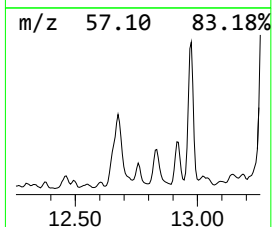
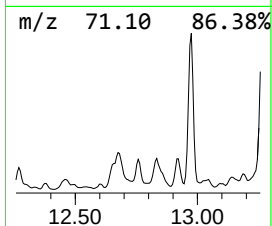
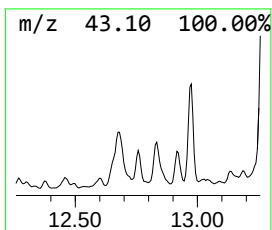
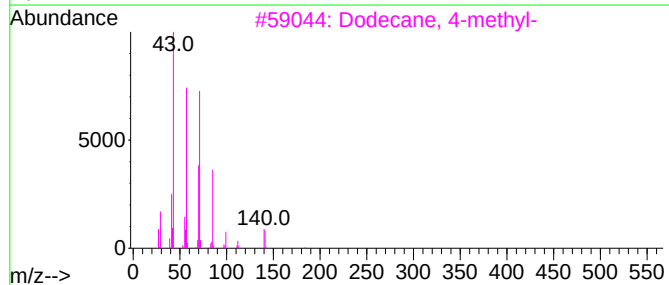
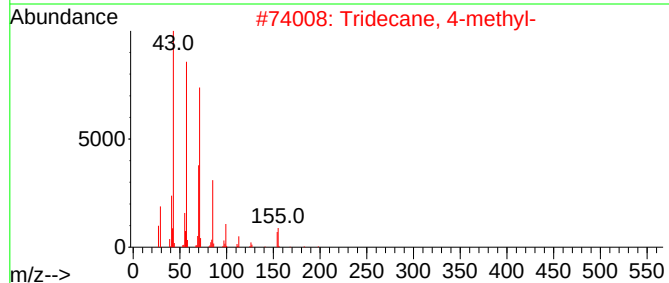
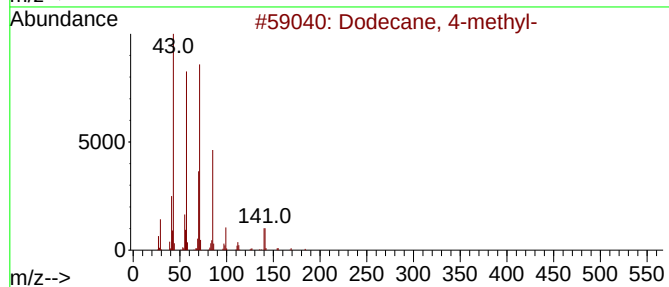
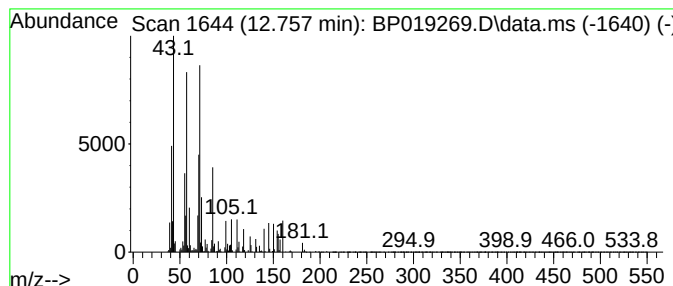
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 (DEL) Alkane: Straight-Chai... Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.757	7.99 ng/ul	726039	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane, 4-methyl-	184	C13H28	006117-97-1	52
2			Tridecane, 4-methyl-	198	C14H30	026730-12-1	52
3			Dodecane, 4-methyl-	184	C13H28	006117-97-1	52
4			Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	50
5			Decane, 5-propyl-	184	C13H28	017312-62-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

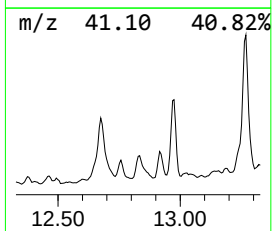
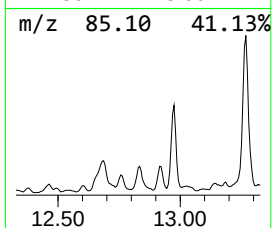
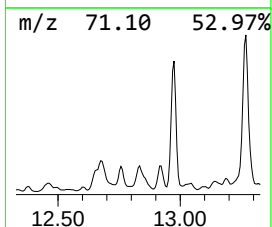
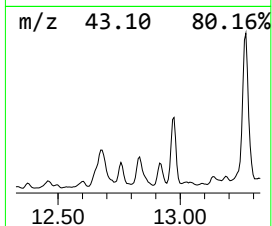
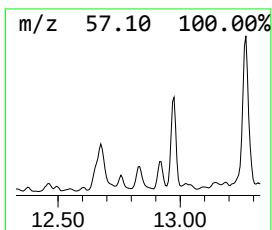
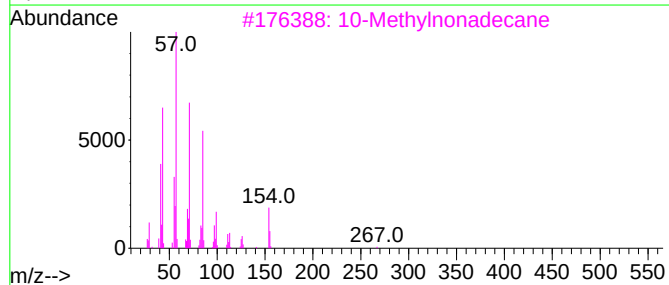
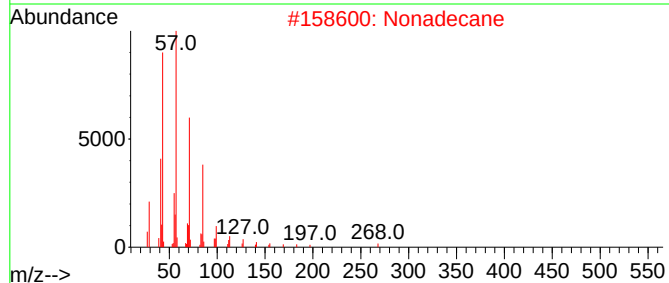
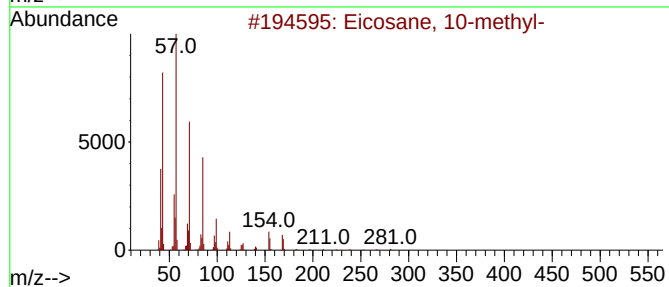
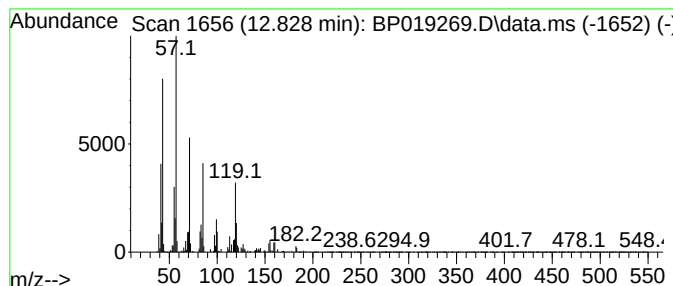
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 32 (DEL) Alkane: Straight-Chai... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.828	12.19 ng/ul	1107480	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Eicosane, 10-methyl-		296	C21H44	054833-23-7	62
2	Nonadecane		268	C19H40	000629-92-5	58
3	10-Methylnonadecane		282	C20H42	056862-62-5	58
4	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	52
5	Tetracosane		338	C24H50	000646-31-1	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

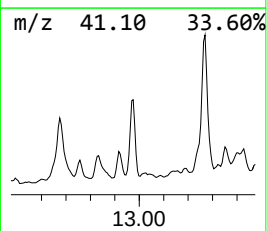
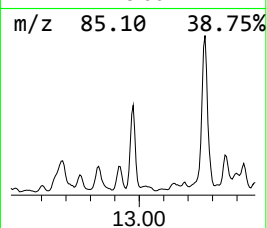
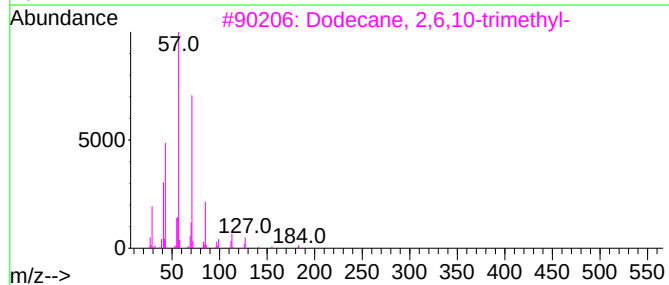
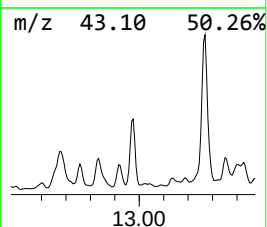
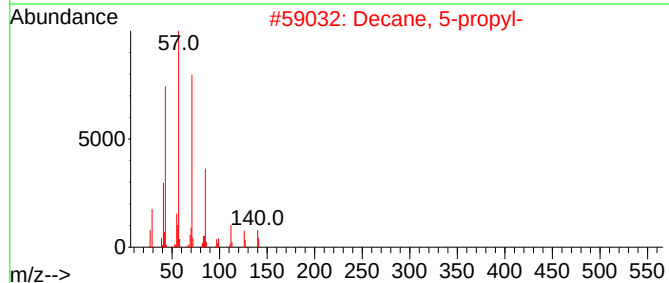
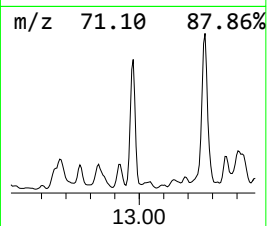
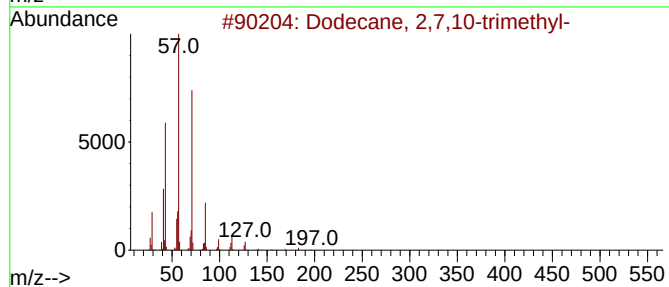
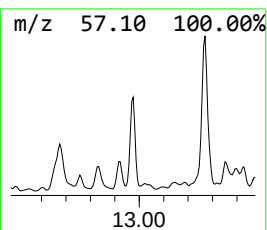
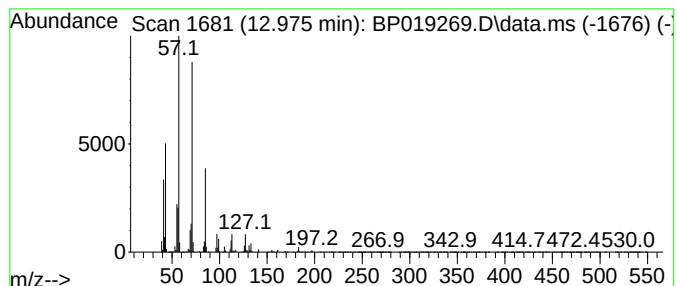
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.975	41.83 ng/ul	3801090	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2,7,10-trimethyl-		212	C15H32	074645-98-0	83
2	Decane, 5-propyl-		184	C13H28	017312-62-8	80
3	Dodecane, 2,6,10-trimethyl-		212	C15H32	003891-98-3	80
4	Nonane, 3,7-dimethyl-		156	C11H24	017302-32-8	76
5	Pentadecane, 7-methyl-		226	C16H34	006165-40-8	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

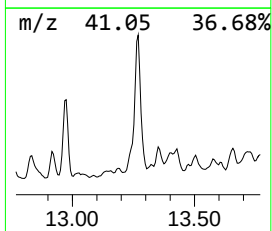
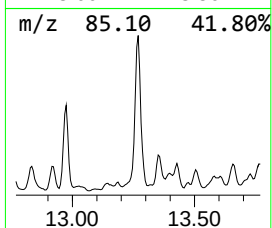
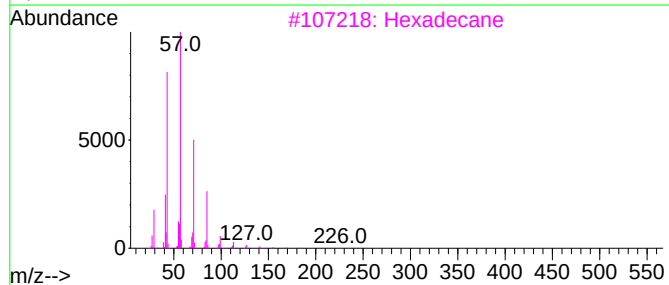
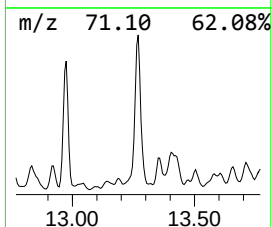
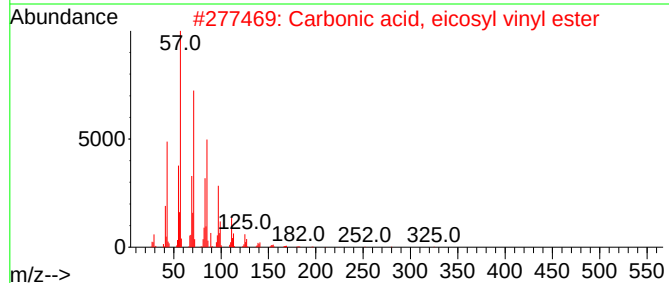
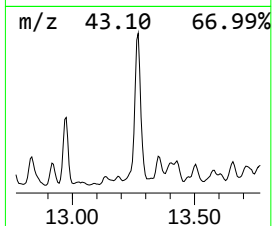
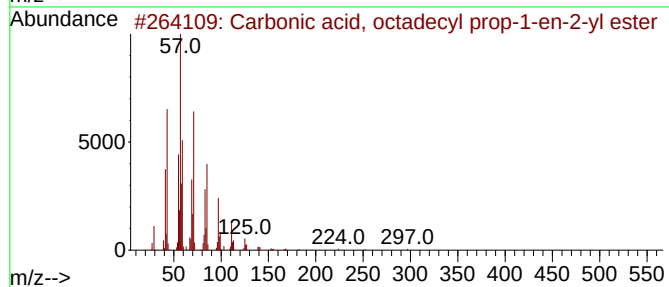
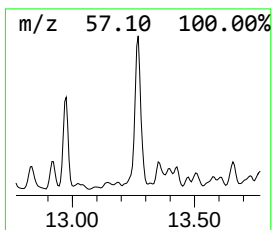
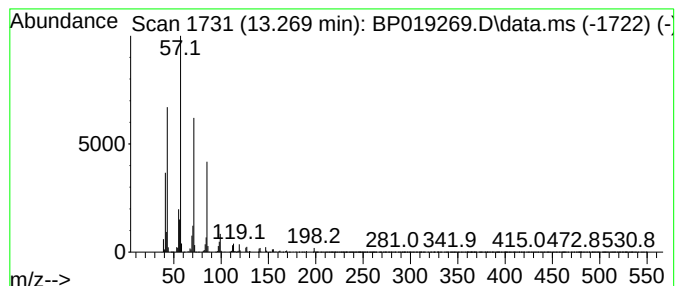
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 36 Carbonic acid, octadecyl pr... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.269	84.25 ng/ul	7654910	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	86
2	Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	80
3	Hexadecane	226	C16H34	000544-76-3	72
4	Hexadecane	226	C16H34	000544-76-3	72
5	Tridecane	184	C13H28	000629-50-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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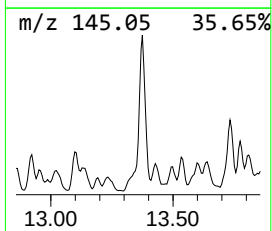
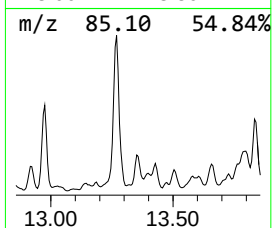
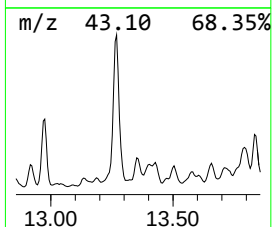
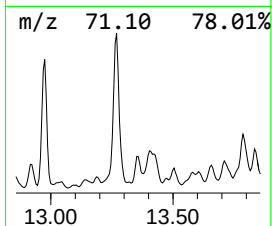
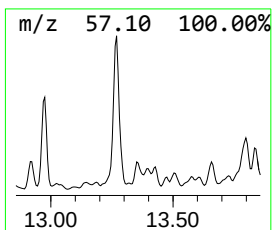
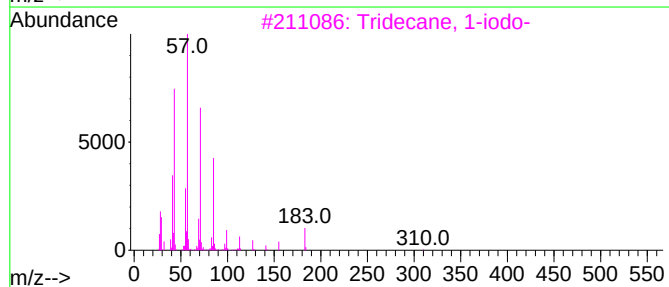
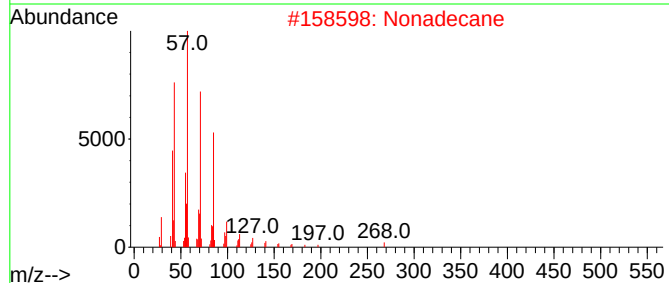
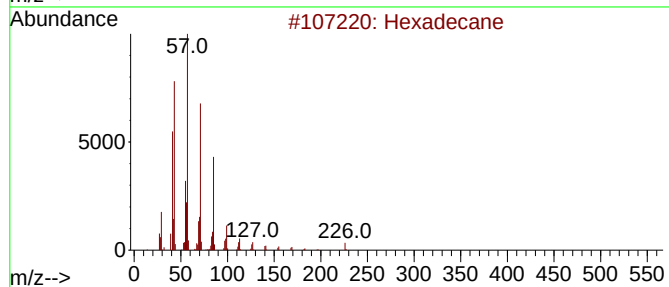
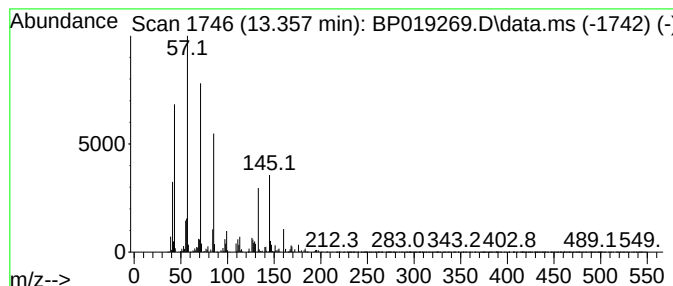
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Quant Title : SVOA CALIBRATION

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

Peak Number 38 (DEL) Alkane: Straight-Chai... Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.357	20.80 ng/ul	1889720	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	58
2			Nonadecane	268	C19H40	000629-92-5	58
3			Tridecane, 1-iodo-	310	C13H27I	035599-77-0	58
4			Nonadecane	268	C19H40	000629-92-5	52
5			Tridecane	184	C13H28	000629-50-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

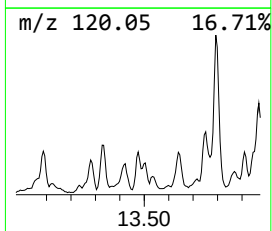
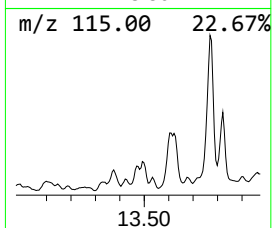
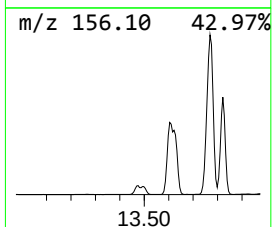
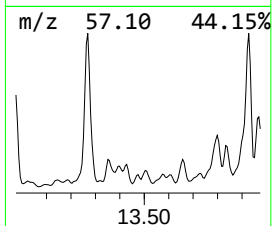
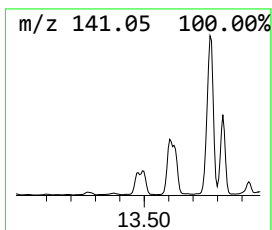
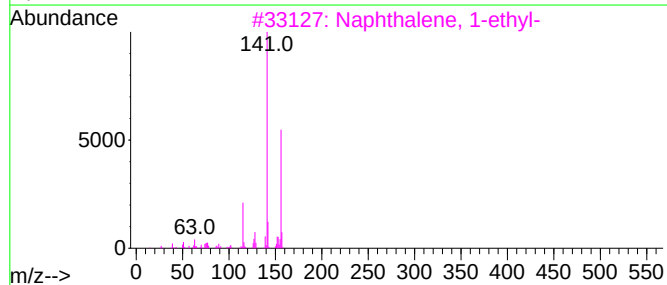
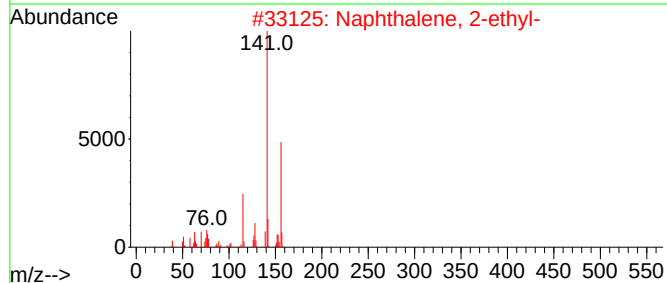
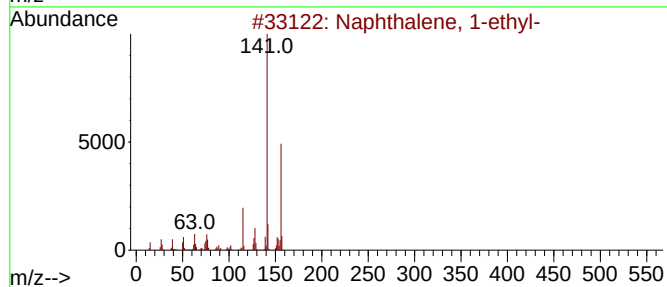
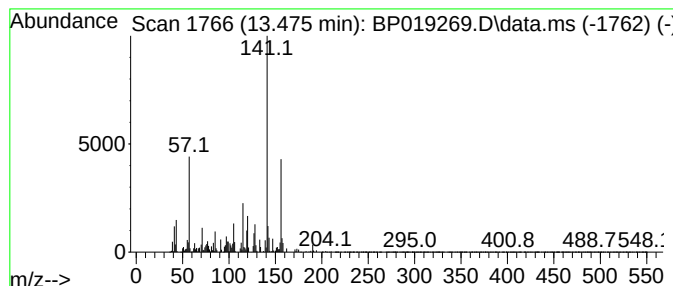
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 Naphthalene, 1-ethyl- Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.475	7.62 ng/ul	691934	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	90
2			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	90
3			Naphthalene, 1-ethyl-	156	C12H12	001127-76-0	89
4			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	81
5			Naphthalene, 2-ethyl-	156	C12H12	000939-27-5	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

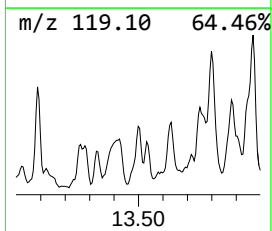
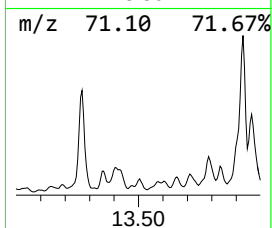
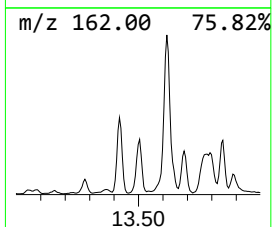
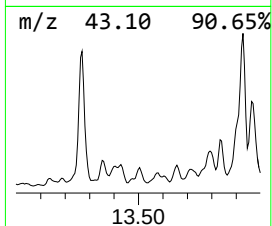
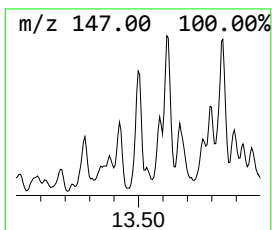
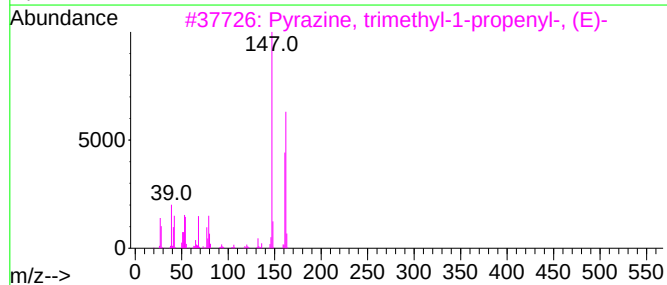
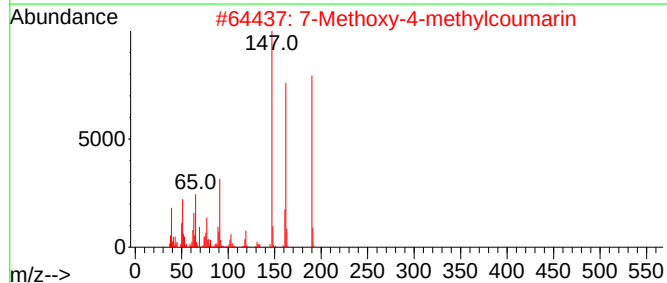
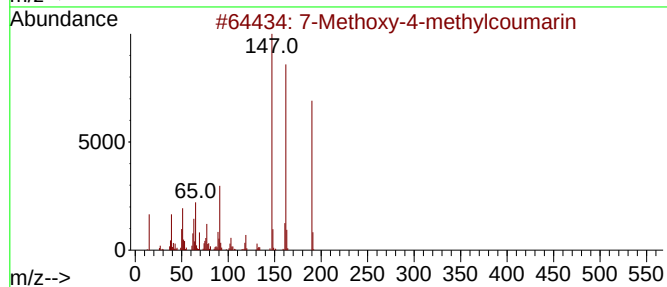
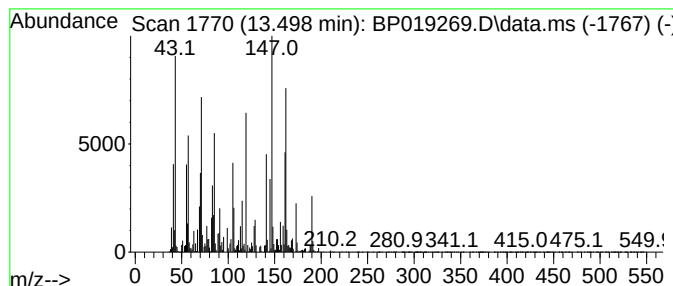
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 7-Methoxy-4-methylcoumarin Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.498	15.51 ng/ul	1409390	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7-Methoxy-4-methylcoumarin	190	C11H10O3	002555-28-4	50
2	7-Methoxy-4-methylcoumarin	190	C11H10O3	002555-28-4	45
3	Pyrazine, trimethyl-1-propenyl-,...	162	C10H14N2	055138-79-9	43
4	Benzene, hexamethyl-	162	C12H18	000087-85-4	38
5	Benzene, hexamethyl-	162	C12H18	000087-85-4	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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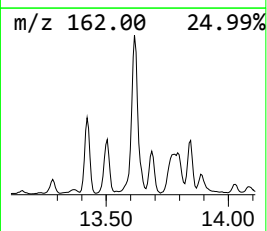
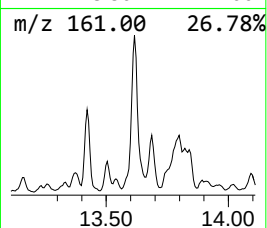
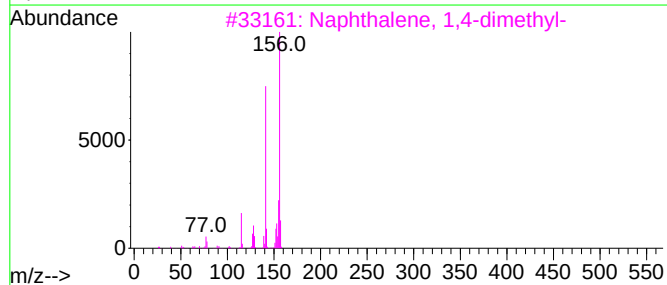
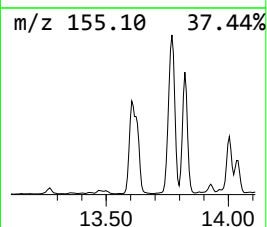
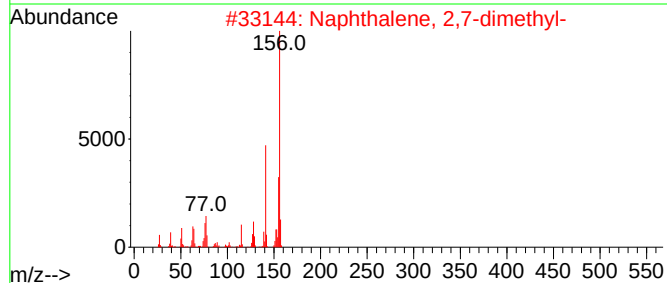
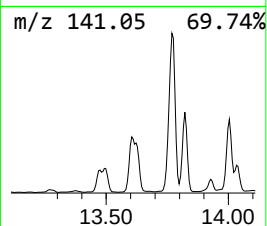
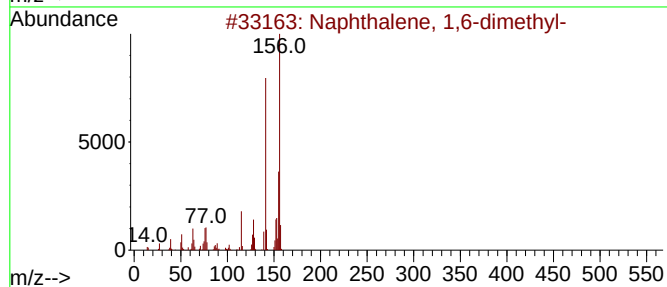
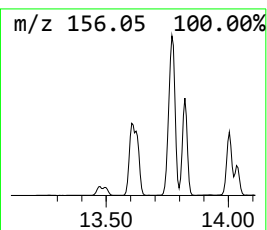
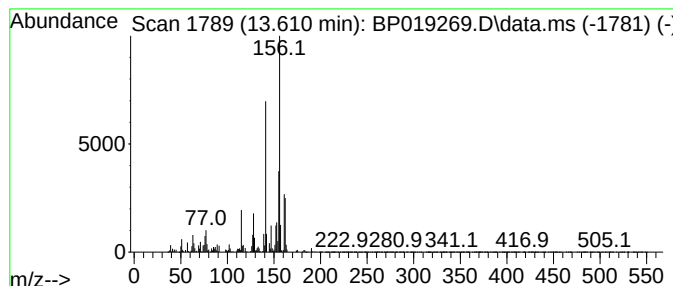
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 Naphthalene, 1,6-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.610	80.46 ng/ul	7310510	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	93
5			Naphthalene, 1,5-dimethyl-	156	C12H12	000571-61-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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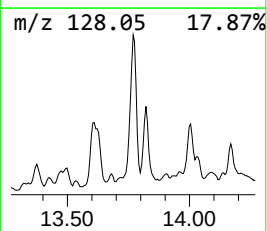
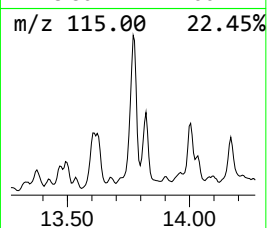
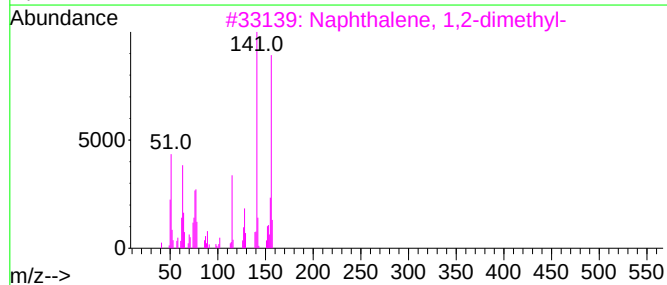
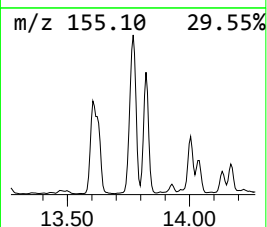
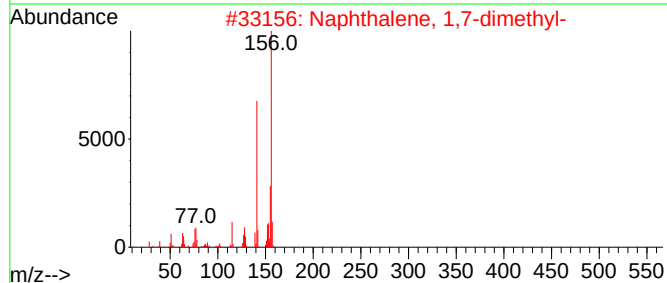
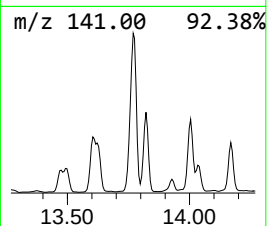
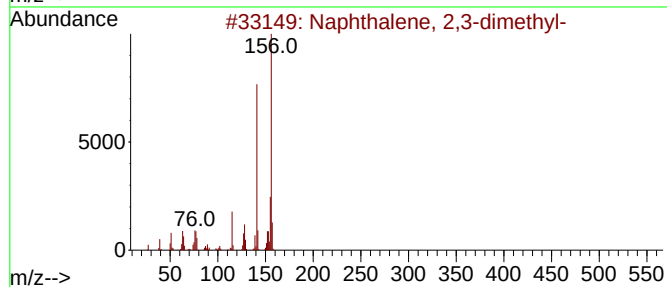
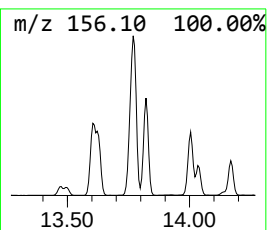
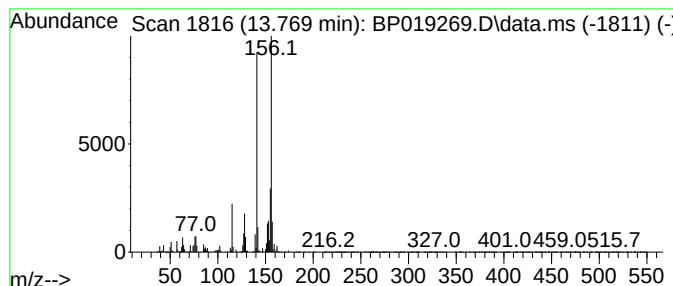
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.769	93.02 ng/ul	8452020	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	97
3			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
5			Naphthalene, 2,6-dimethyl-	156	C12H12	000581-42-0	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

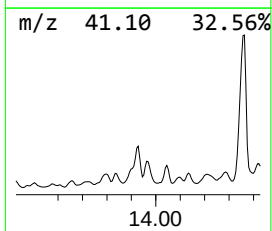
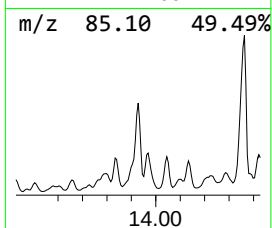
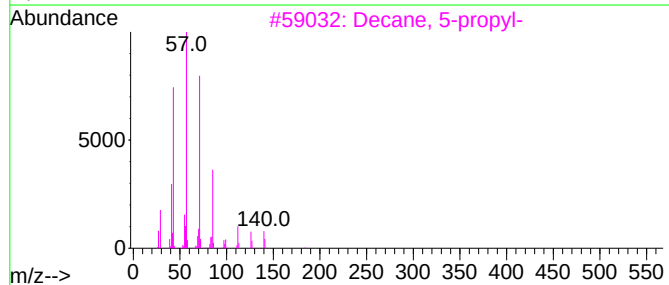
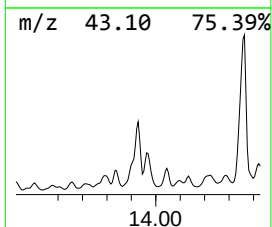
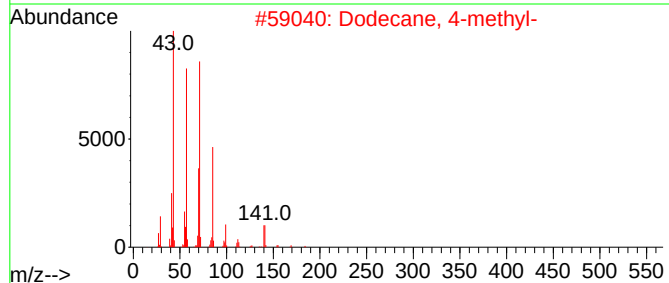
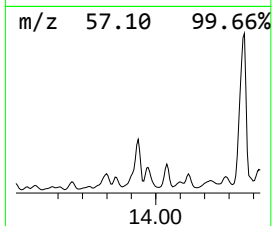
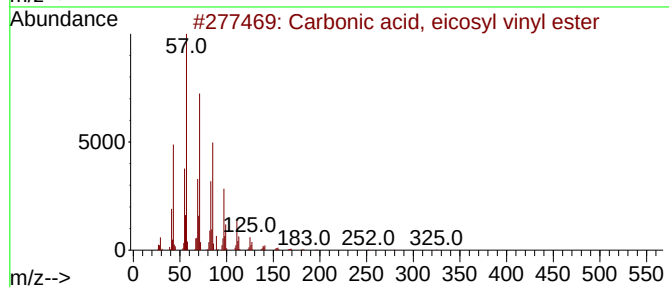
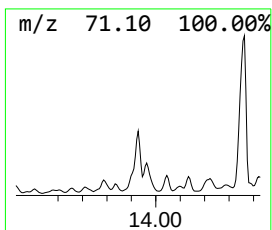
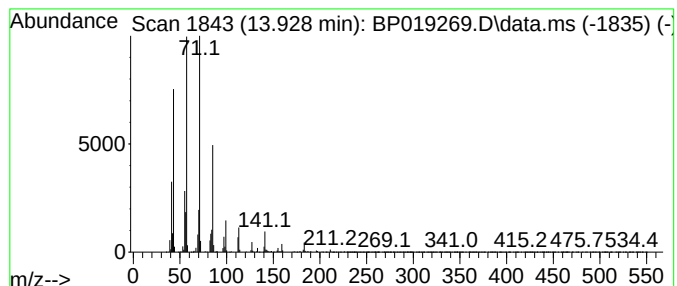
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 Carbonic acid, eicosyl vinyl... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.928	87.91 ng/ul	7987870	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Carbonic acid, eicosyl vinyl ester	368	C23H44O3	1000382-54-3	86
2			Dodecane, 4-methyl-	184	C13H28	006117-97-1	81
3			Decane, 5-propyl-	184	C13H28	017312-62-8	81
4			Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	80
5			Tridecane, 5-propyl-	226	C16H34	055045-11-9	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

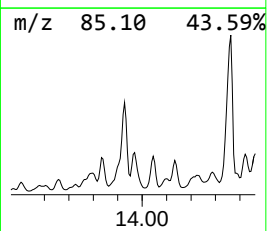
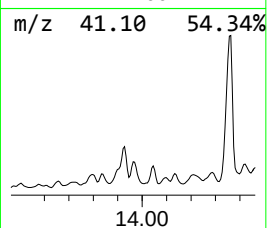
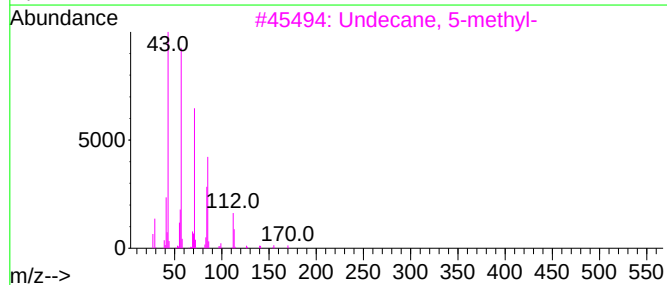
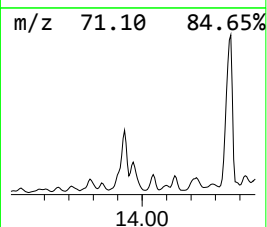
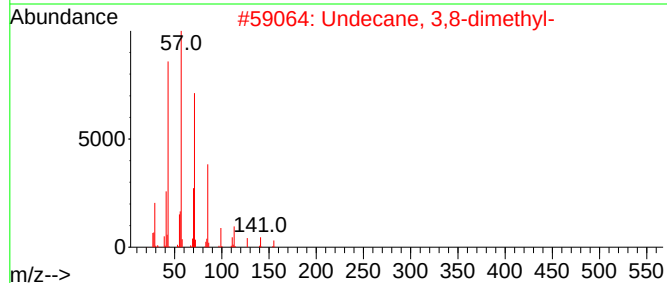
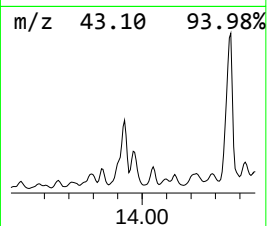
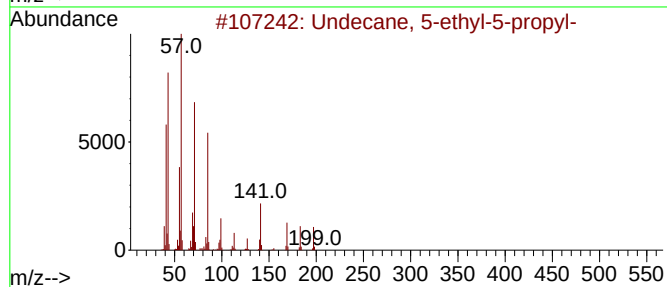
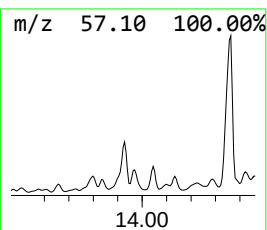
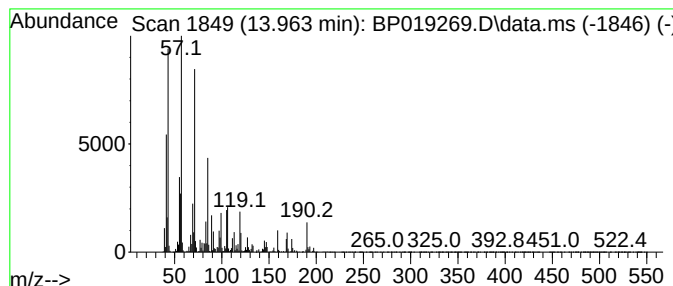
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.963	38.42 ng/ul	3491300	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	50
2	Undecane, 3,8-dimethyl-	184	C13H28	017301-30-3	49
3	Undecane, 5-methyl-	170	C12H26	001632-70-8	49
4	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	49
5	Hexatriacontane	507	C36H74	000630-06-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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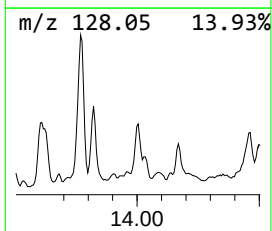
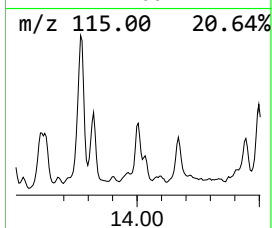
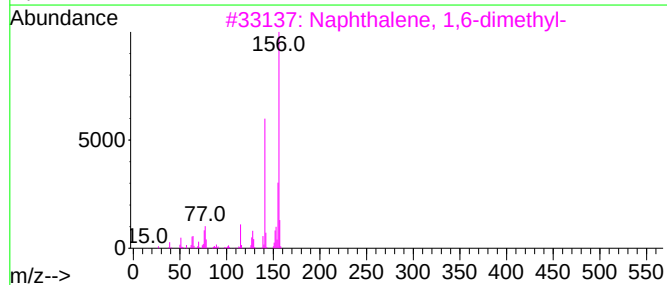
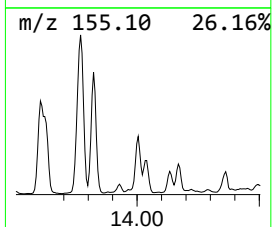
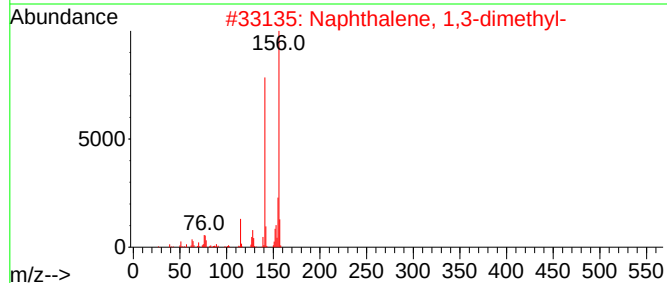
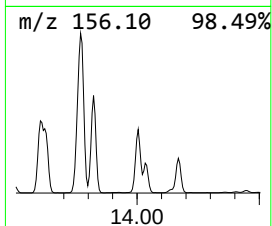
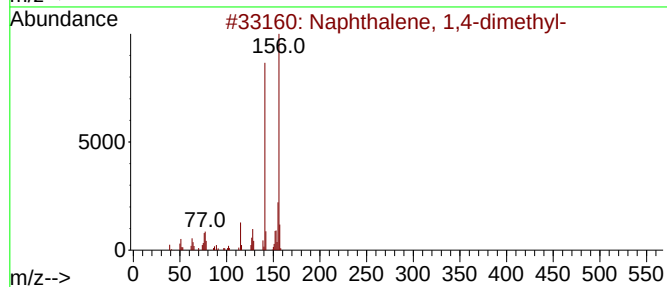
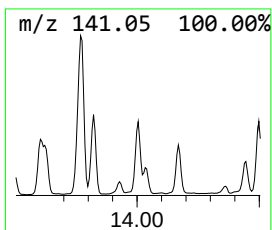
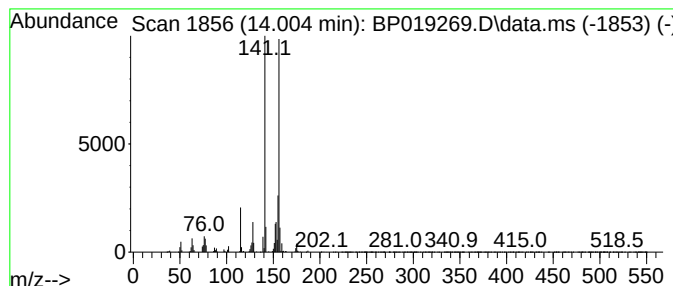
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Naphthalene, 1,4-dimethyl- Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.004	16.18 ng/ul	1469980	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
2			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	94
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

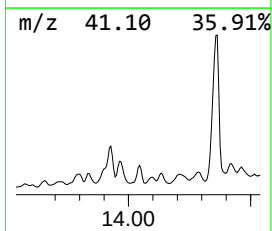
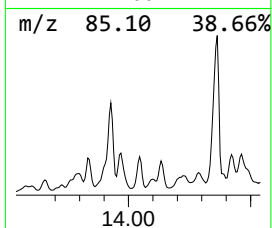
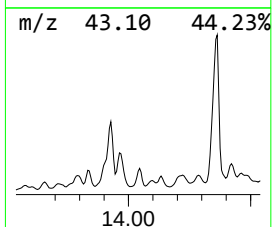
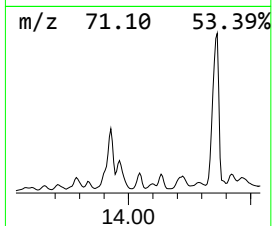
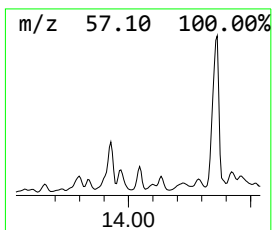
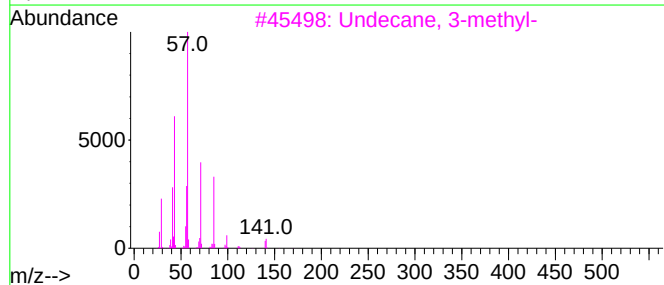
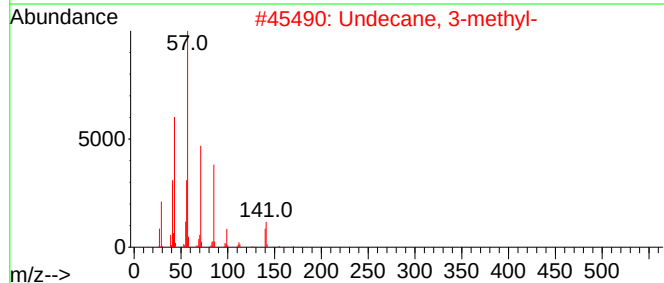
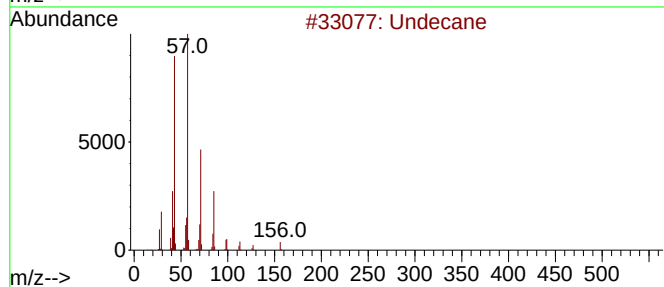
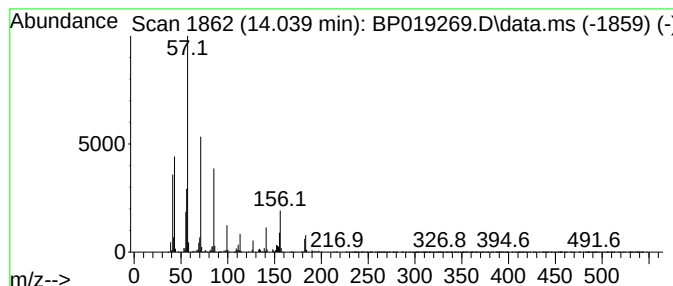
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 46 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
14.040	33.96 ng/ul	3085840	Acenaphthene-d10		14.445	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Undecane		156	C11H24	001120-21-4	74
2	Undecane, 3-methyl-		170	C12H26	001002-43-3	64
3	Undecane, 3-methyl-		170	C12H26	001002-43-3	59
4	Borane, diethyl(decyloxy)-		226	C14H31BO	1000152-34-3	53
5	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

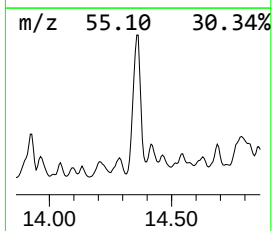
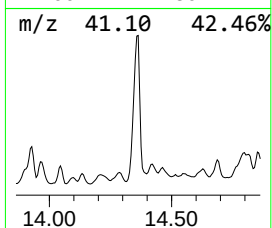
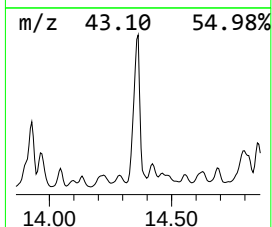
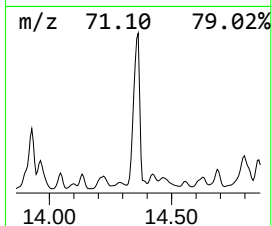
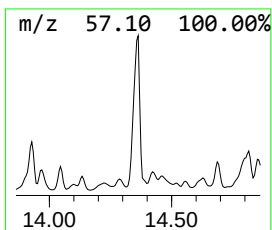
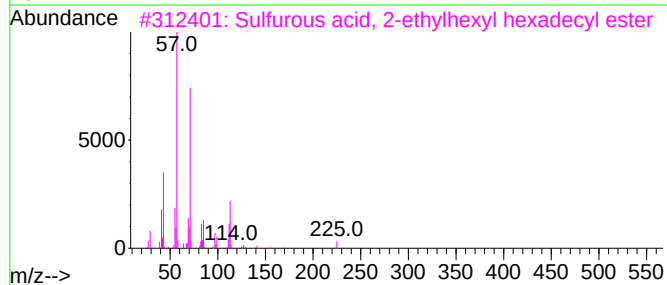
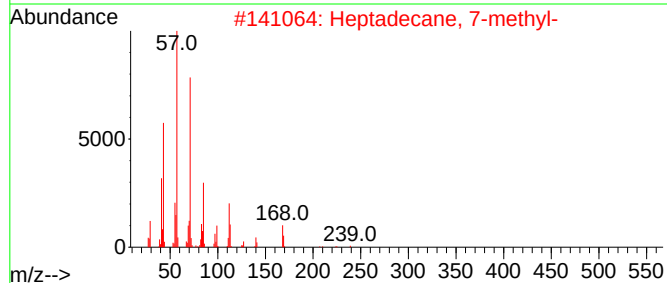
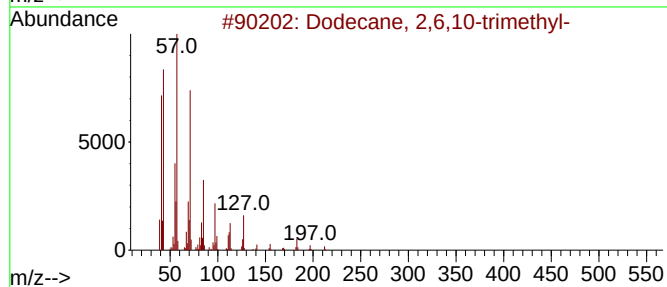
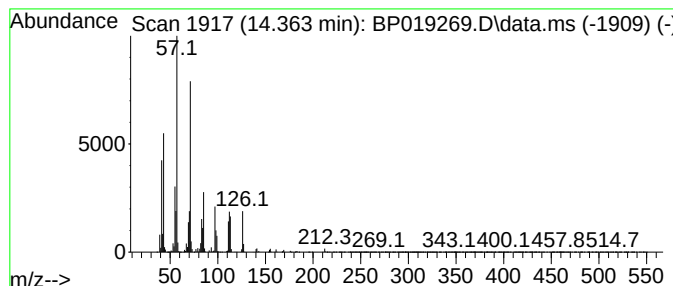
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Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.363	346.87 ng/ul	31516600	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Dodecane, 2,6,10-trimethyl-	212	C15H32	003891-98-3	83
2	Heptadecane, 7-methyl-	254	C18H38	020959-33-5	59
3	Sulfurous acid, 2-ethylhexyl hex...	418	C24H50O3S	1000309-19-9	53
4	Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	53
5	Dodecane, 2,6,11-trimethyl-	212	C15H32	031295-56-4	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

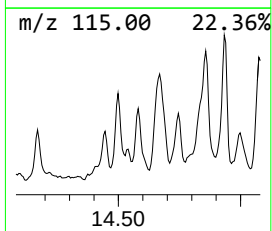
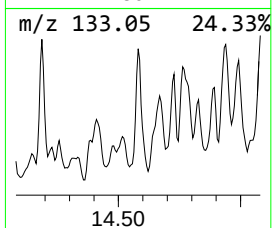
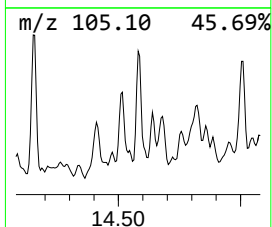
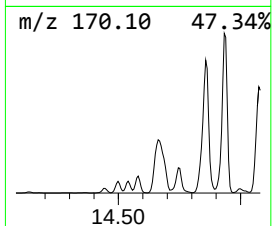
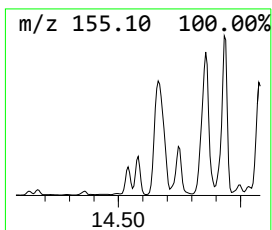
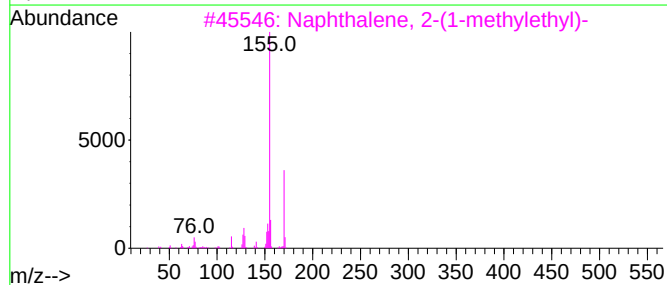
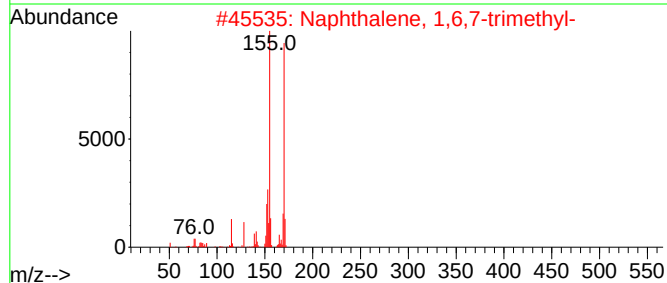
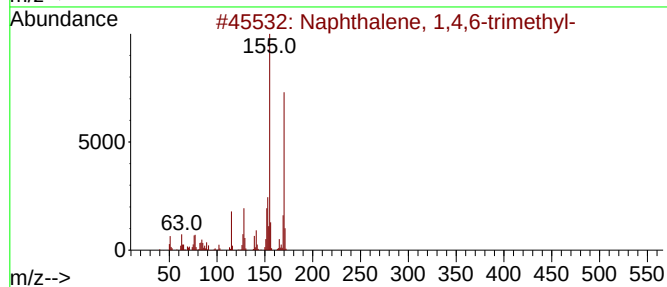
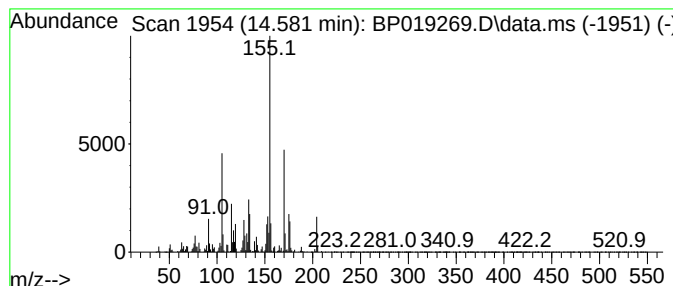
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 Naphthalene, 1,4,6-trimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.581	20.24 ng/ul	1838980	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	90
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	86
3	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	70
4	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	70
5	Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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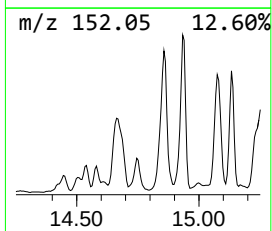
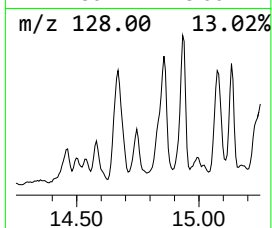
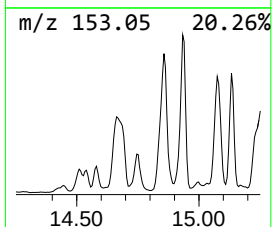
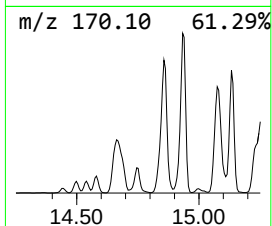
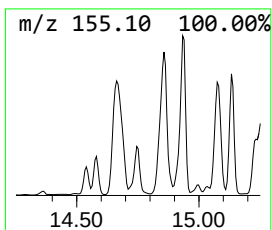
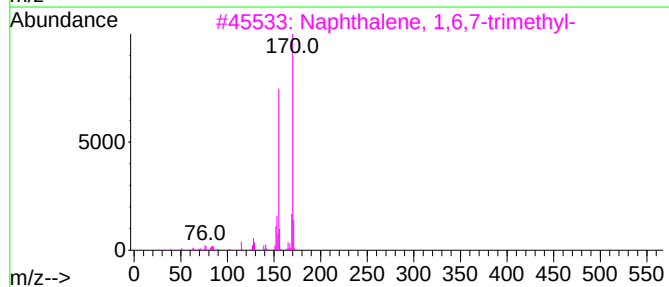
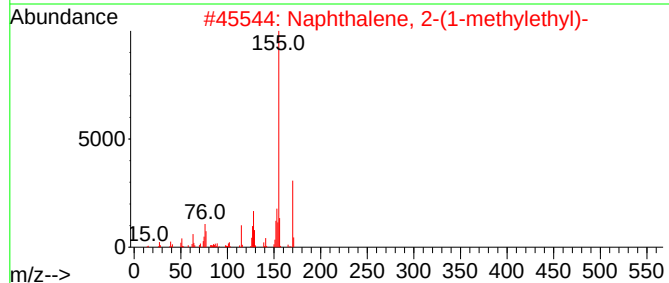
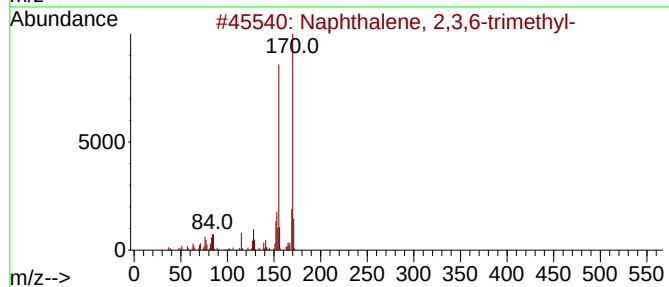
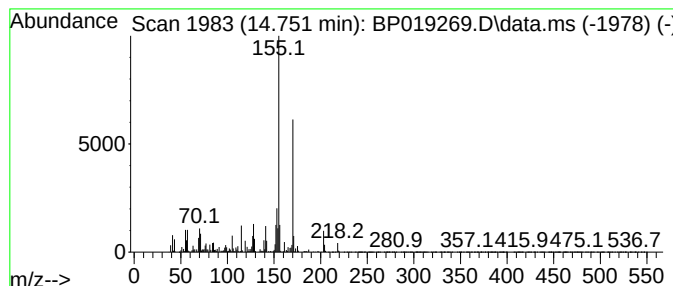
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 Naphthalene, 2,3,6-trimethyl- Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.751	42.25 ng/ul	3838470	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93
2			Naphthalene, 2-(1-methylethyl)-	170	C13H14	002027-17-0	93
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
4			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	92
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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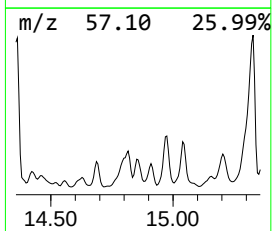
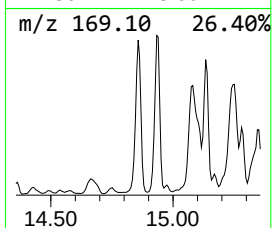
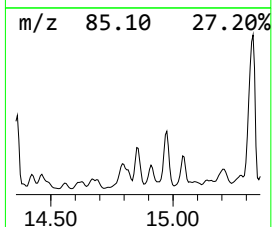
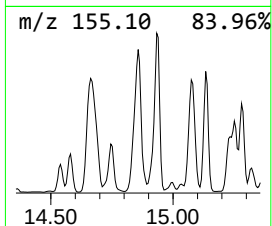
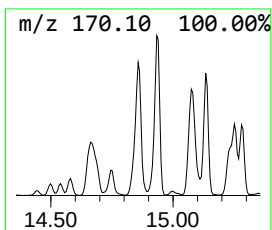
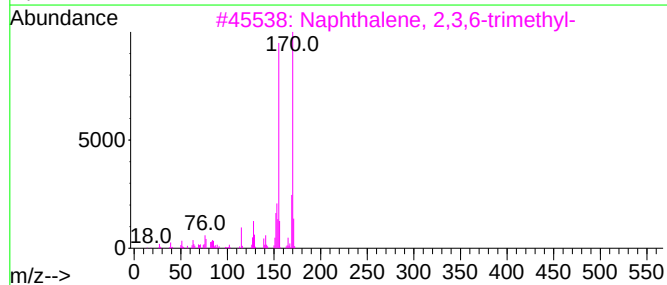
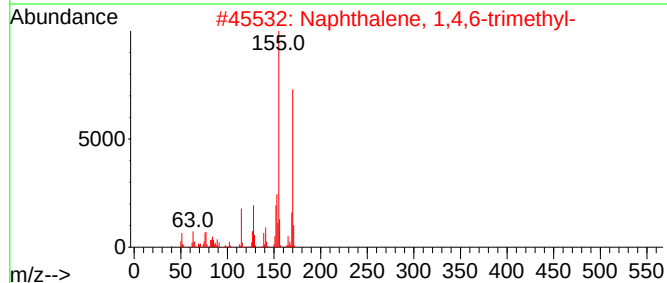
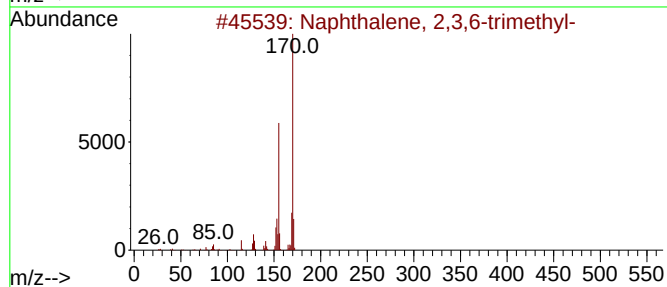
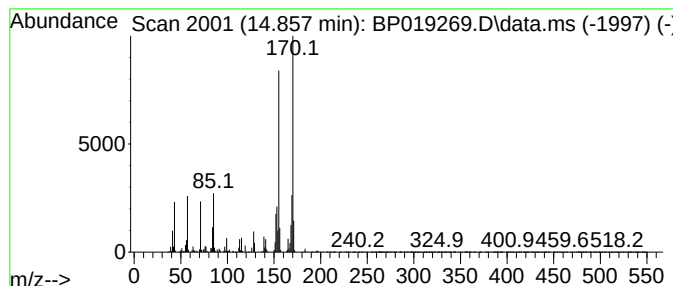
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Quant Title : SVOA CALIBRATION

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

Peak Number 50 Naphthalene, 1,6,7-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.857	136.17 ng/ul	12372800	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	95
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	95
3			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
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Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

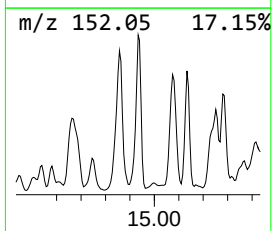
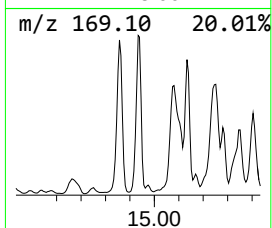
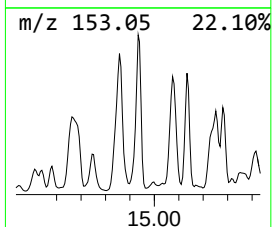
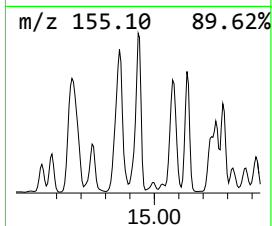
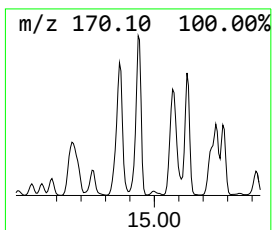
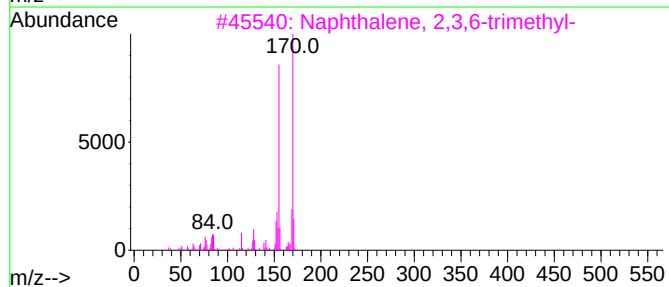
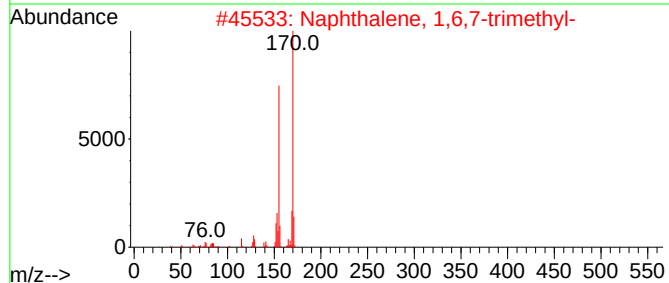
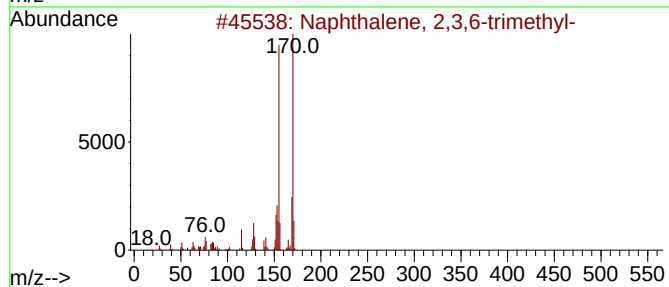
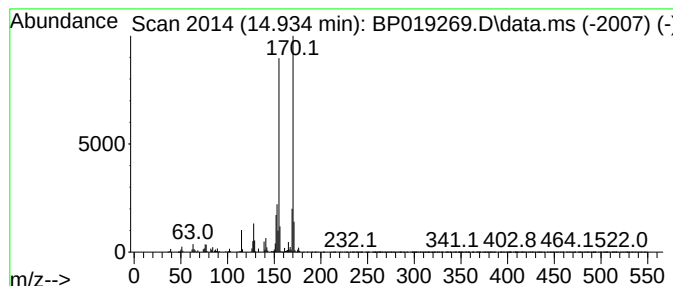
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BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 unknown-05 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.934	125.45 ng/ul	11398600	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	98
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	97
3	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
4	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	96
5	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

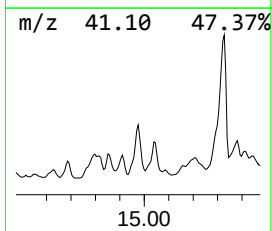
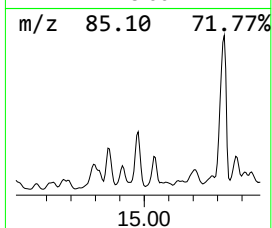
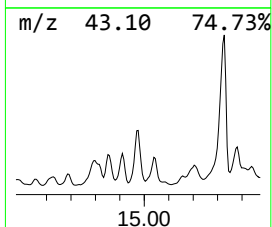
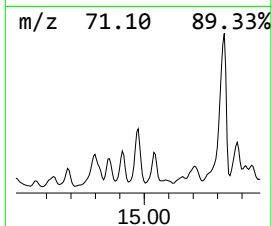
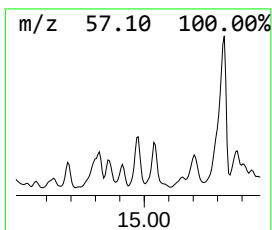
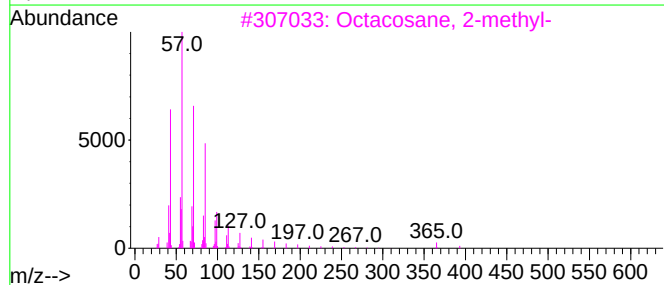
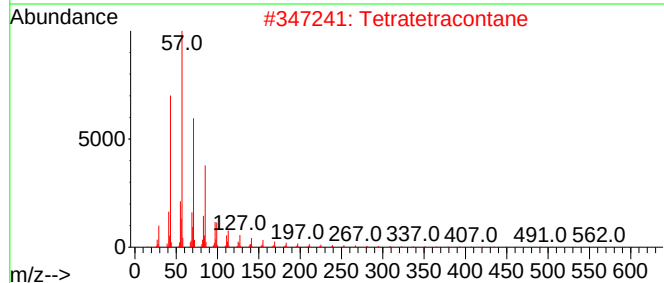
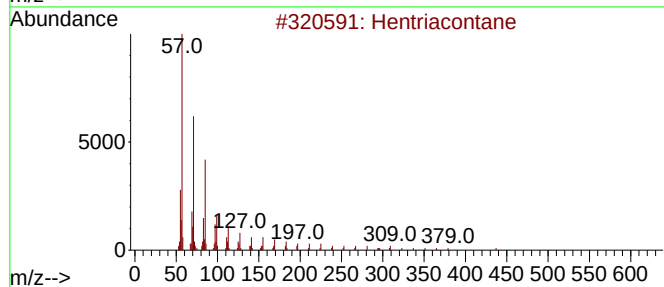
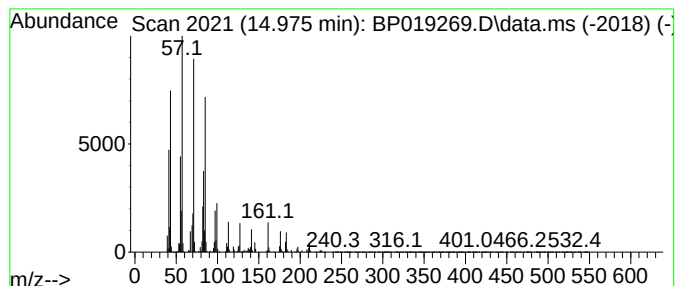
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 52 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.975	80.59 ng/ul	7322830	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hentriacontane	437	C31H64	000630-04-6	86
2	Tetratetracontane	619	C44H90	007098-22-8	80
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	80
4	Hexacosane	366	C26H54	000630-01-3	80
5	Dotriacontane, 1-iodo-	576	C32H65I	1000406-32-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

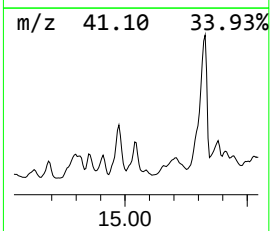
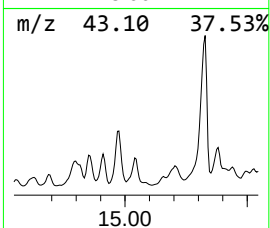
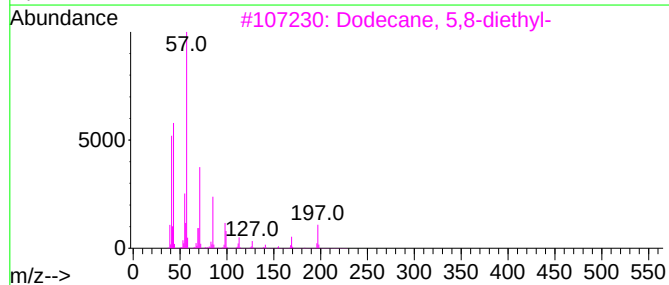
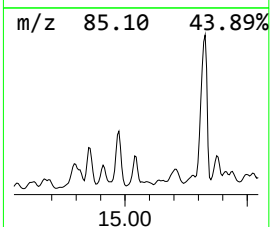
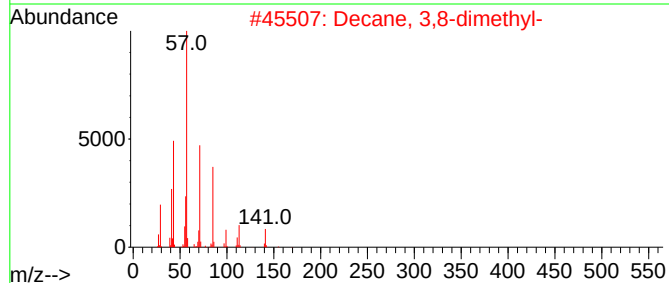
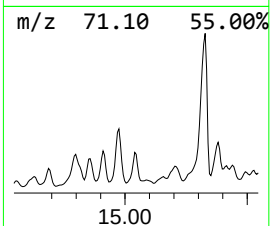
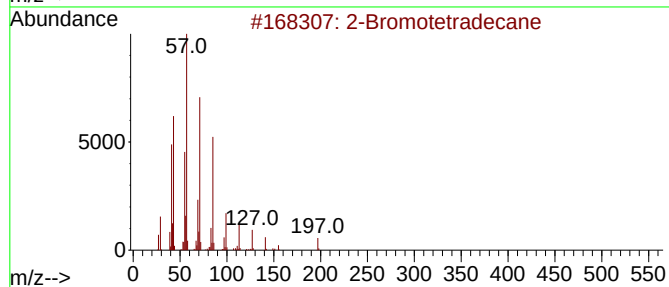
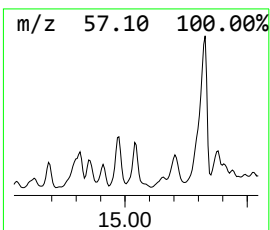
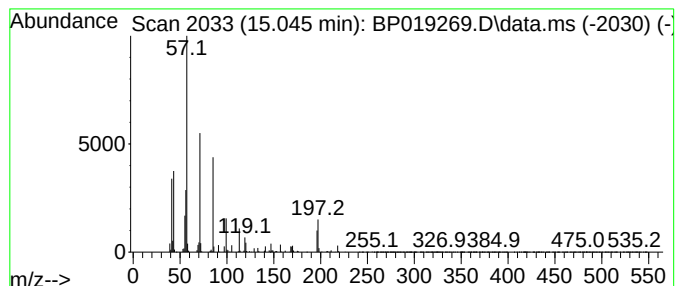
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 53 2-Bromotetradecane Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.045	29.43 ng/ul	2674160	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Bromotetradecane	276	C14H29Br	074036-95-6	72
2	Decane, 3,8-dimethyl-	170	C12H26	017312-55-9	72
3	Dodecane, 5,8-diethyl-	226	C16H34	024251-86-3	64
4	Undecane, 3-methyl-	170	C12H26	001002-43-3	64
5	Undecane, 5-ethyl-5-propyl-	226	C16H34	002755-07-9	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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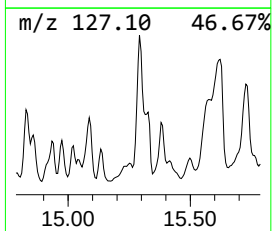
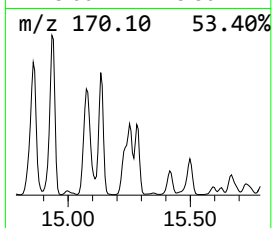
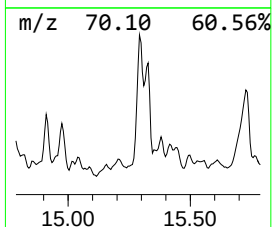
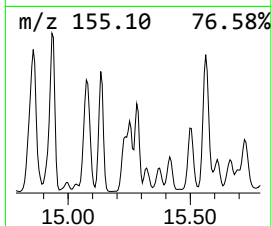
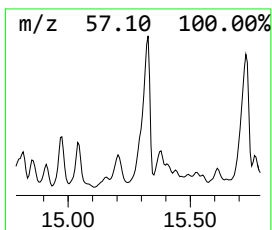
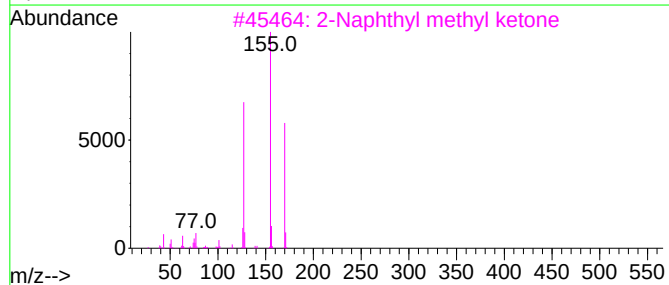
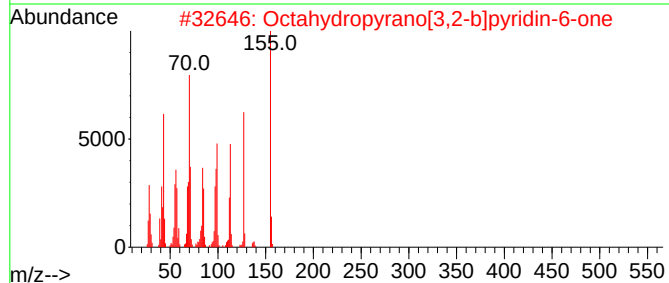
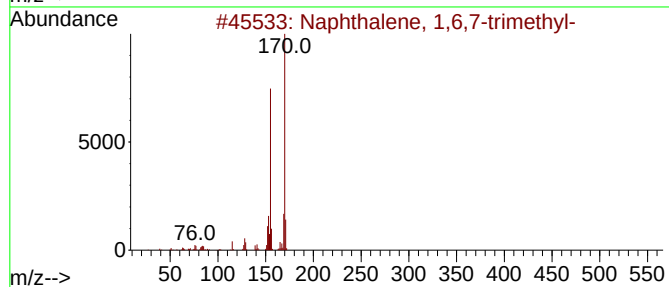
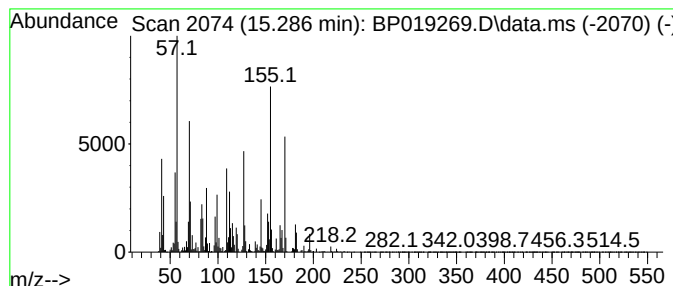
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Quant Title : SVOA CALIBRATION

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

Peak Number 54 unknown-06 Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.286	85.47 ng/ul	7765700	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	49
2			Octahydropyrano[3,2-b]pyridin-6-one	155	C8H13NO2	1000193-43-8	49
3			2-Naphthyl methyl ketone	170	C12H10O	000093-08-3	43
4			Cis-4-methyl-cis-2-thiabicyclo[4...	170	C10H18S	1000491-53-0	43
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	42



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

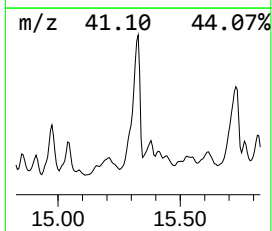
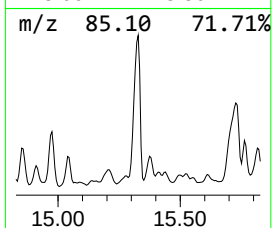
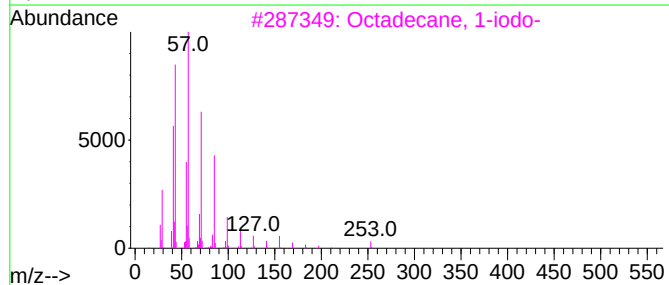
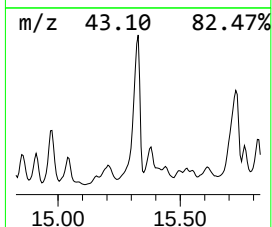
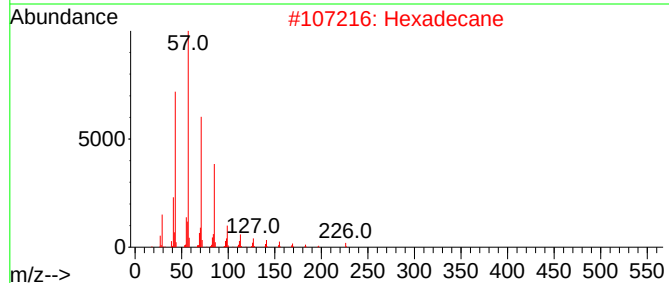
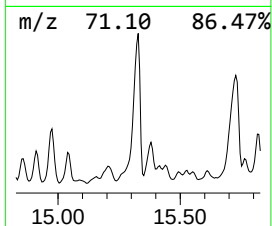
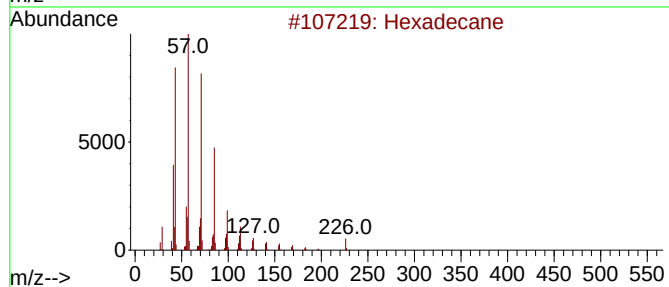
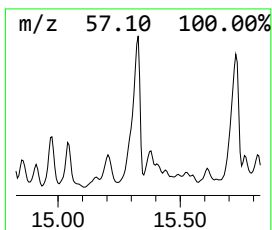
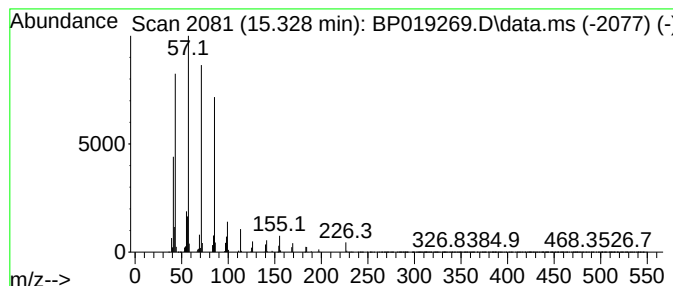
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.328	177.83 ng/ul	16157900	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	95
2	Hexadecane	226	C16H34	000544-76-3	91
3	Octadecane, 1-iodo-	380	C18H37I	000629-93-6	90
4	Hexadecane	226	C16H34	000544-76-3	86
5	Octacosane	394	C28H58	000630-02-4	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

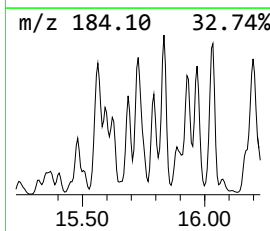
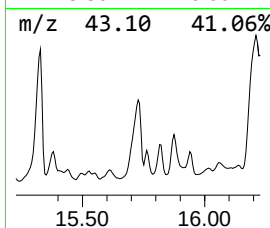
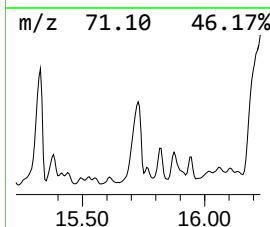
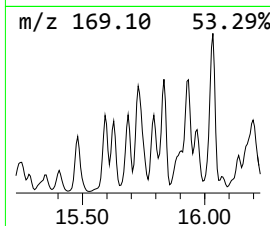
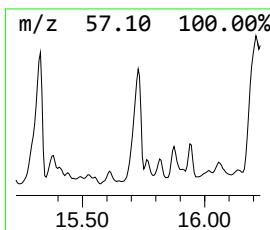
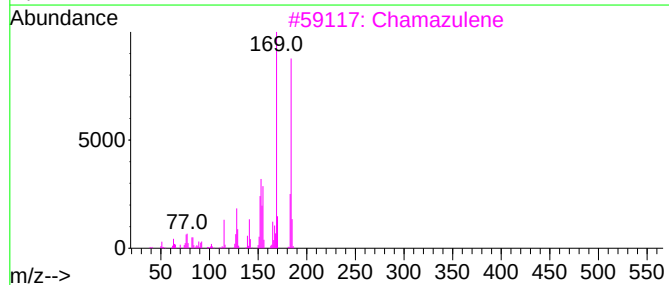
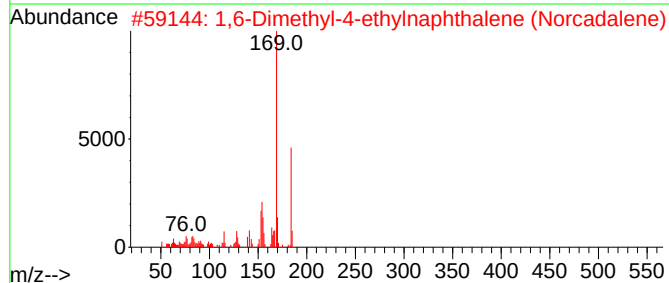
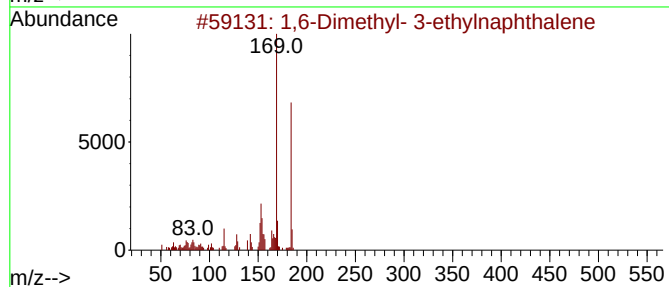
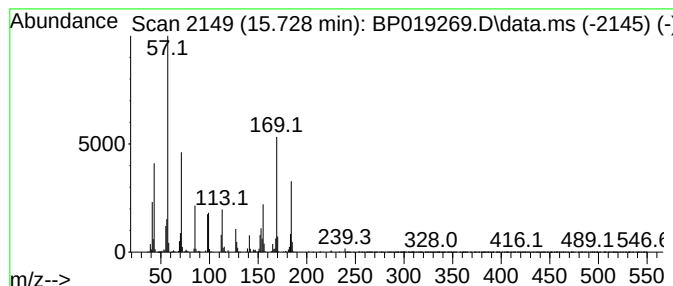
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Quant Title : SVOA CALIBRATION

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

Peak Number 56 unknown-07 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.728	171.72 ng/ul	15602900	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,6-Dimethyl- 3-ethylnaphthalene	184	C14H16	1000383-71-8	42
2			1,6-Dimethyl-4-ethylnaphthalene ...	184	C14H16	1000383-71-7	38
3			Chamazulene	184	C14H16	000529-05-5	35
4			1-Naphthohydroxamic acid	187	C11H9NO2	006953-61-3	35
5			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

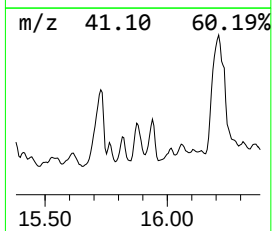
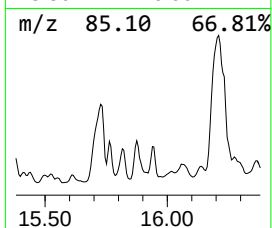
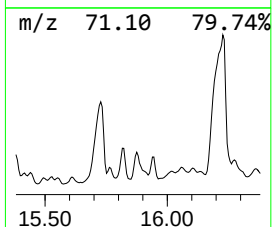
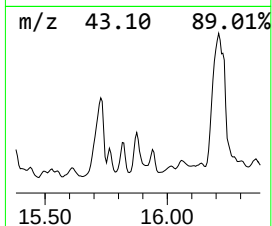
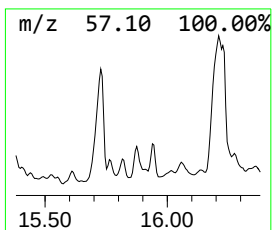
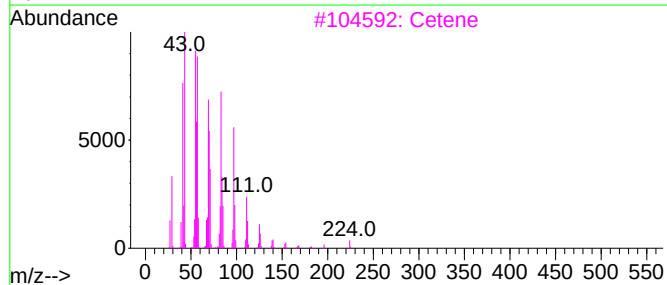
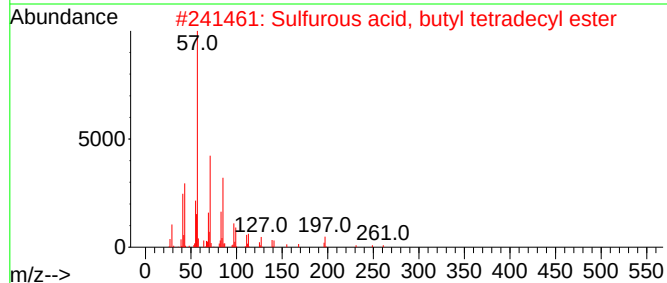
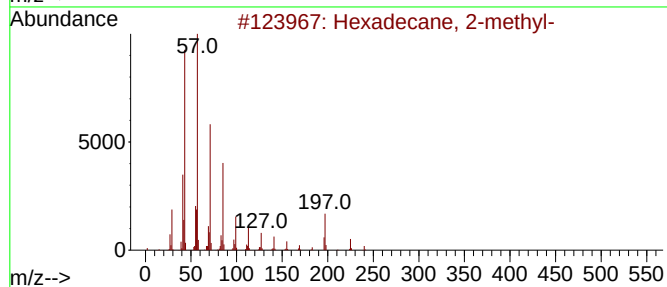
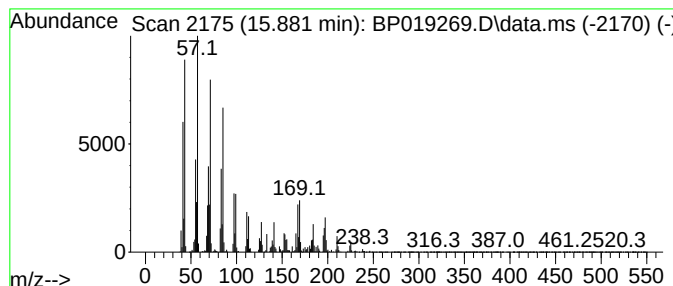
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Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 48

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.881	7.32 ng/ul	10776500	Phenanthrene-d10	17.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	90
2			Sulfurous acid, butyl tetradecyl...	334	C18H38O3S	1010309-18-1	81
3			Cetene	224	C16H32	000629-73-2	56
4			Cetene	224	C16H32	000629-73-2	56
5			1-Heptadecene	238	C17H34	006765-39-5	56



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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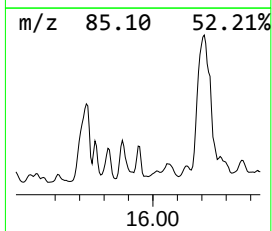
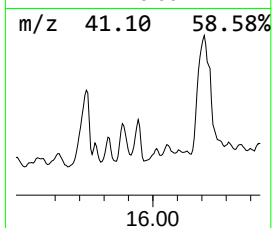
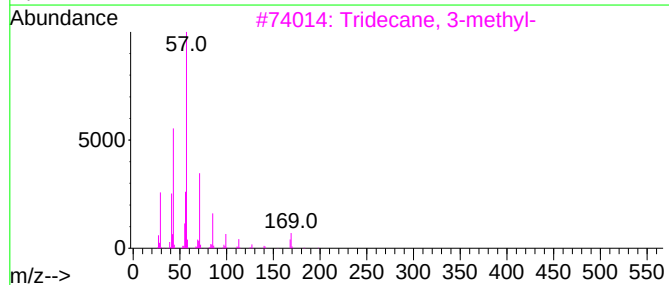
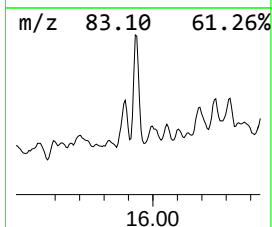
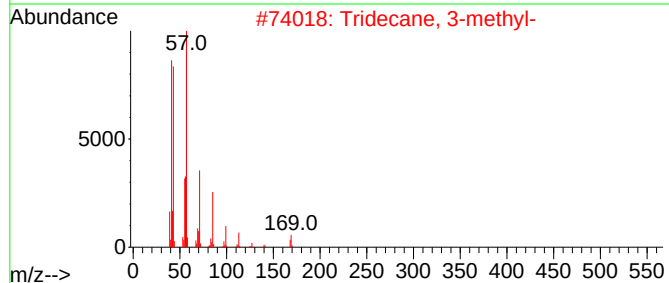
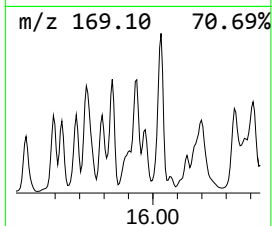
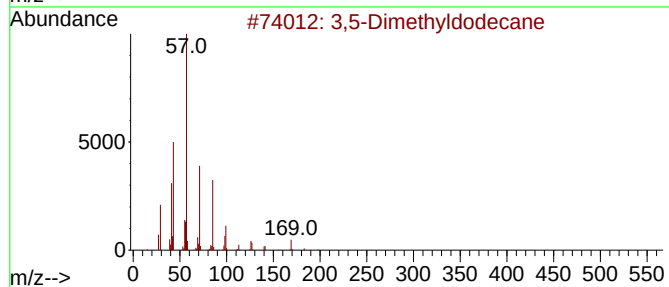
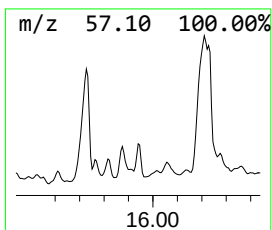
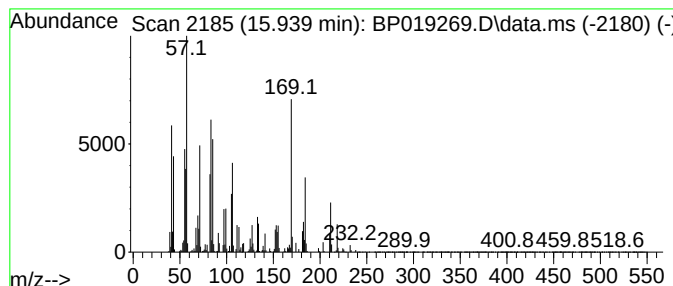
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 (DEL) Alkane: Straight-Chai... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.939	6.27 ng/ul	9234510	Phenanthrene-d10	17.275

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		3,5-Dimethyldodecane	198	C14H30	107770-99-0	42
2		Tridecane, 3-methyl-	198	C14H30	006418-41-3	41
3		Tridecane, 3-methyl-	198	C14H30	006418-41-3	30
4		Pentadecane, 7-(bromomethyl)-	304	C16H33Br	052997-43-0	27
5		Dodecane, 3-methyl-	184	C13H28	017312-57-1	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

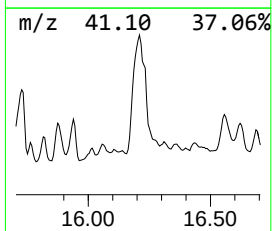
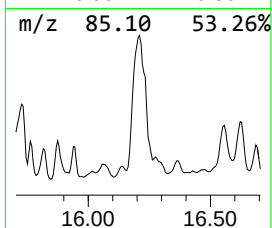
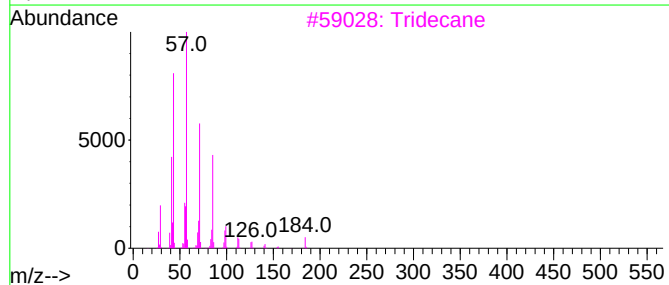
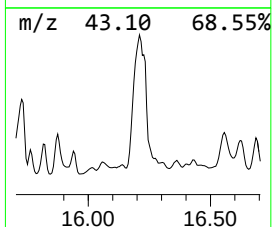
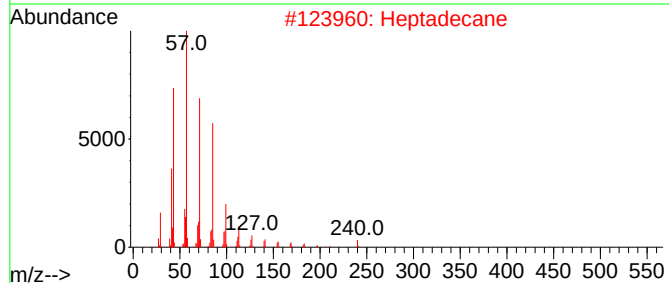
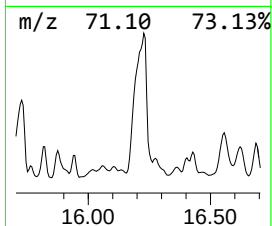
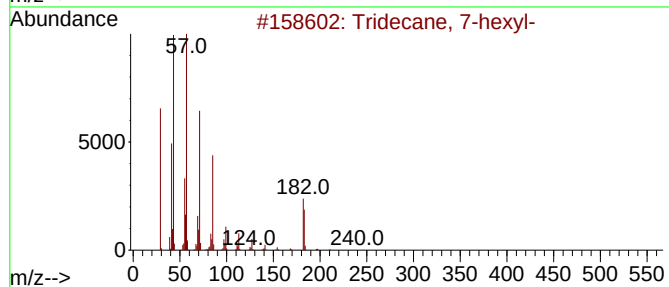
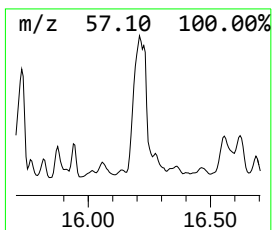
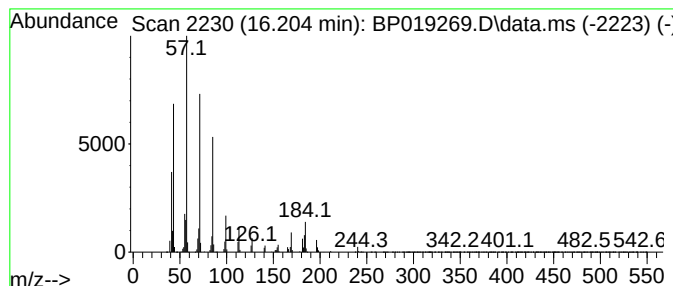
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Quant Title : SVOA CALIBRATION

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

Peak Number 59 (DEL) Alkane: Straight-Chai... Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.204	20.77 ng/ul	30566700	Phenanthrene-d10	17.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 7-hexyl-	268	C19H40	007225-66-3	91
2			Heptadecane	240	C17H36	000629-78-7	87
3			Tridecane	184	C13H28	000629-50-5	83
4			Octadecane	254	C18H38	000593-45-3	80
5			Sulfurous acid, 2-propyl tetra...	320	C17H36O3S	1010309-12-5	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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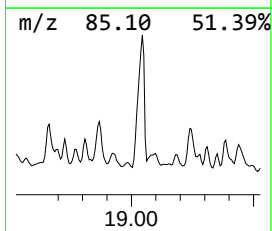
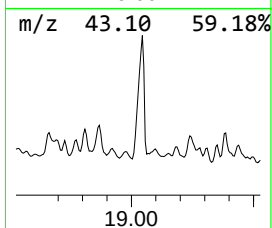
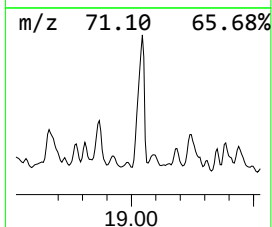
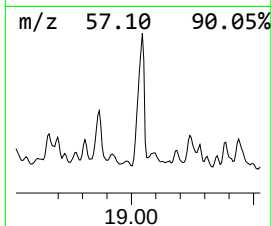
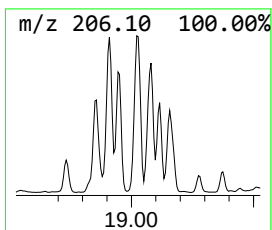
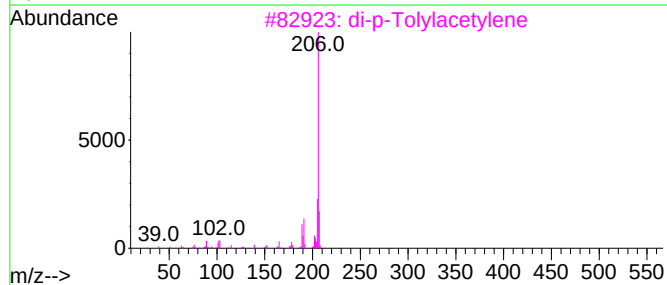
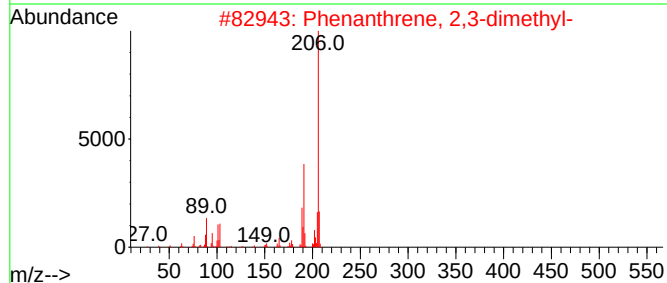
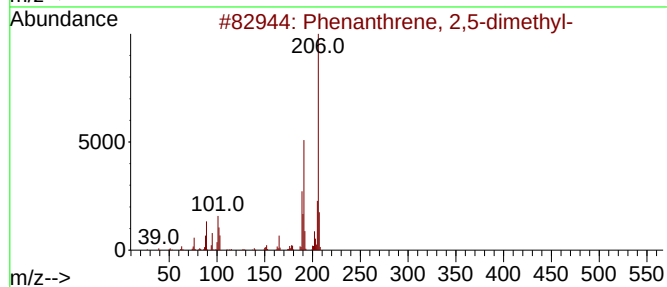
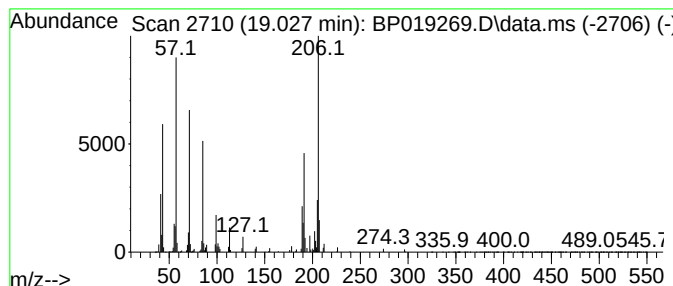
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 Phenanthrene, 2,5-dimethyl- Concentration Rank 41

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.027	9.23 ng/ul	13587900	Phenanthrene-d10	17.275

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenanthrene, 2,5-dimethyl-	206	C16H14	003674-66-6	92
2			Phenanthrene, 2,3-dimethyl-	206	C16H14	003674-65-5	60
3			di-p-Tolylacetylene	206	C16H14	002789-88-0	60
4			Pentadecane	212	C15H32	000629-62-9	53
5			Phenanthrene, 2,7-dimethyl-	206	C16H14	001576-69-8	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCF81

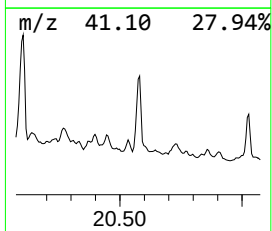
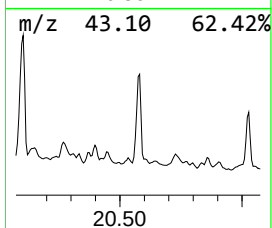
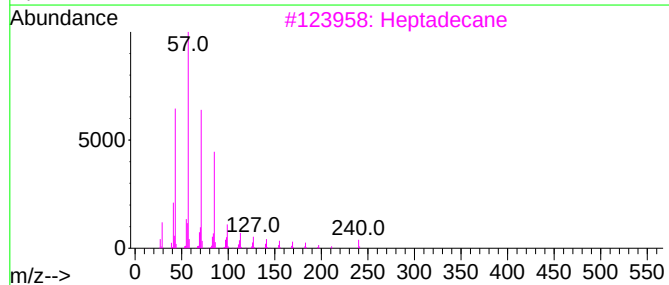
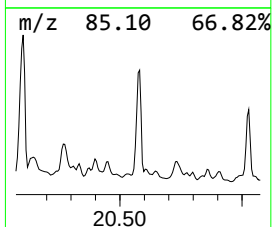
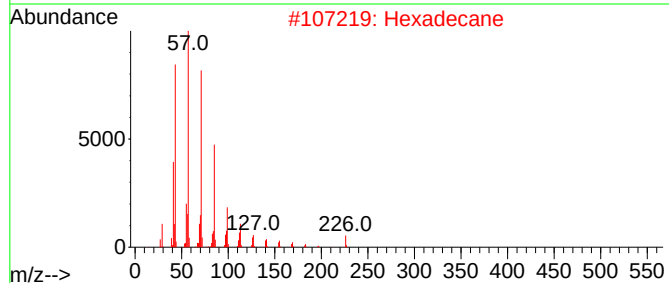
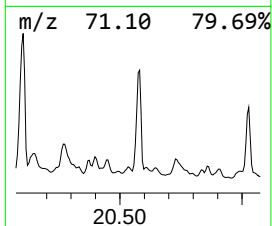
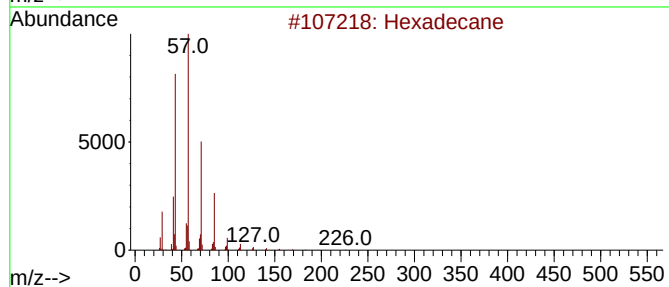
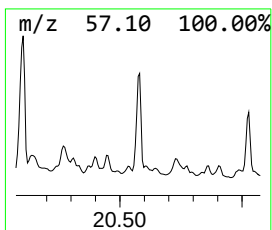
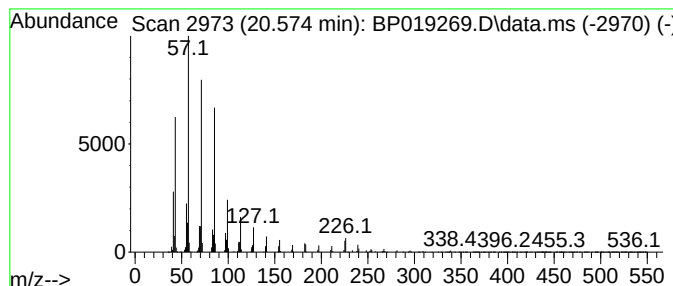
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.574	28.29 ng/ul	13331900	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexadecane	226	C16H34	000544-76-3	96
2			Hexadecane	226	C16H34	000544-76-3	95
3			Heptadecane	240	C17H36	000629-78-7	95
4			Octadecane	254	C18H38	000593-45-3	94
5			Hexadecane, 2-methyl-	240	C17H36	001560-92-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

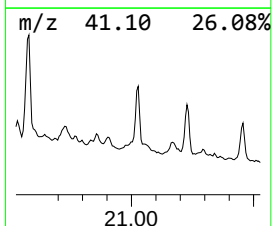
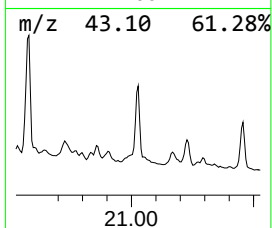
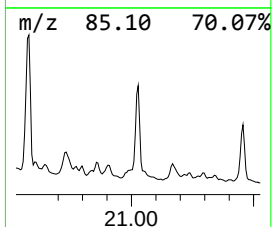
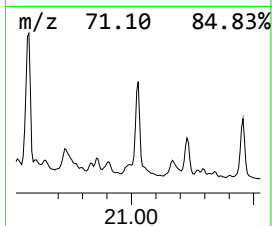
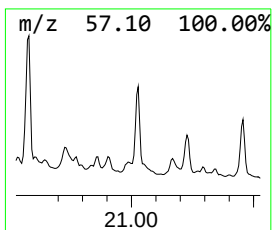
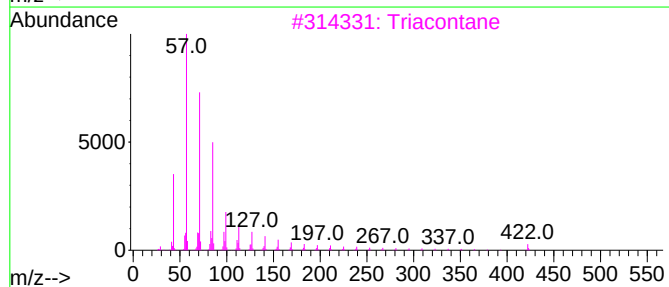
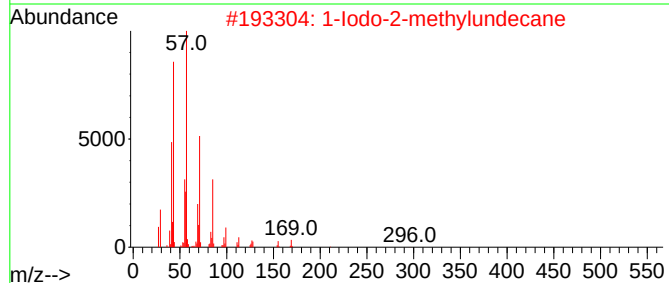
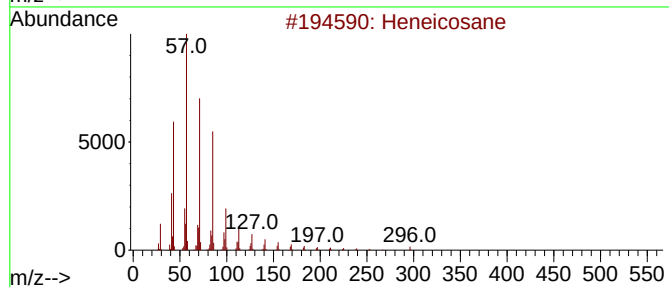
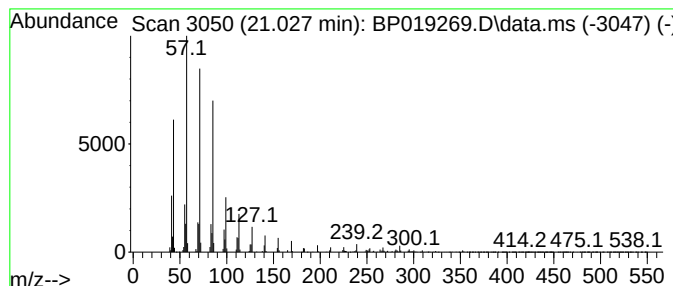
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 62 (DEL) Alkane: Straight-Chai... Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.027	13.64 ng/ul	6428840	Chrysene-d12		21.333	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane		296	C21H44	000629-94-7	99
2	1-Iodo-2-methylundecane		296	C12H25I	073105-67-6	91
3	Triacontane		422	C30H62	000638-68-6	91
4	Octacosane		394	C28H58	000630-02-4	91
5	Pentacosane		352	C25H52	000629-99-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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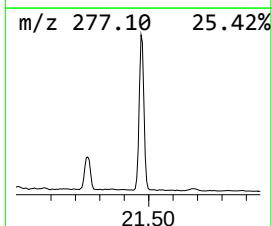
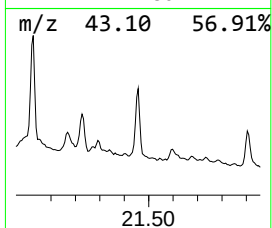
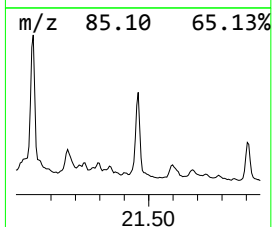
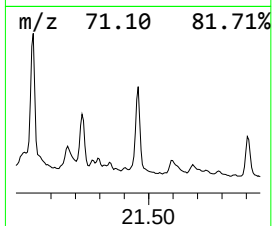
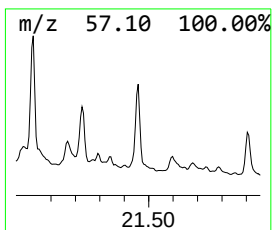
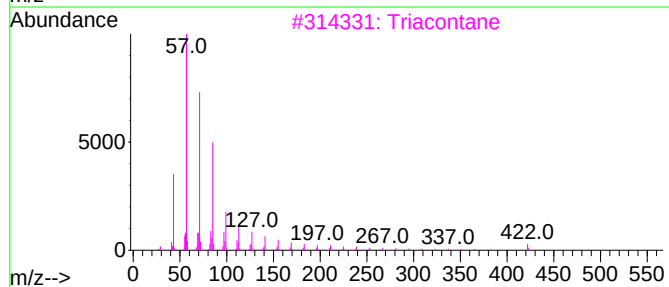
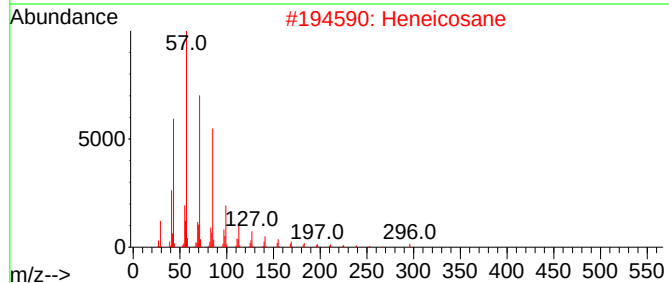
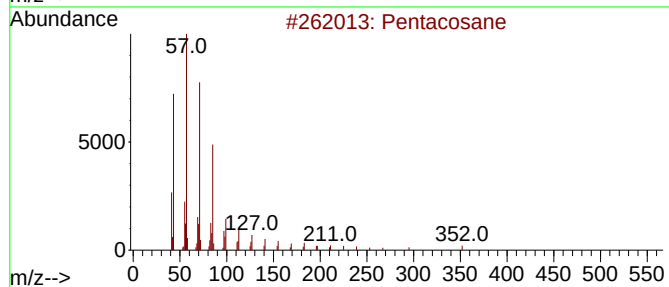
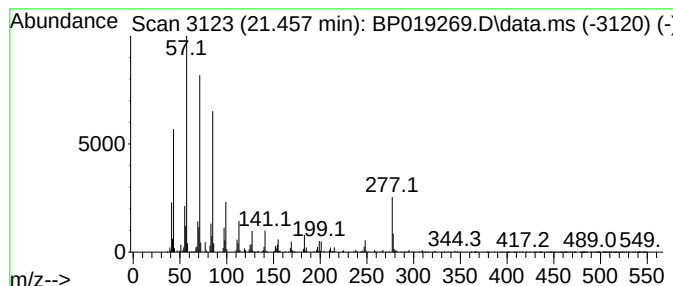
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 (DEL) Alkane: Straight-Chai... Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.457	14.76 ng/ul	6953890	Chrysene-d12	21.333

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacosane	352	C25H52	000629-99-2	90
2			Heneicosane	296	C21H44	000629-94-7	87
3			Triacontane	422	C30H62	000638-68-6	87
4			Hexacosane, 1-iodo-	492	C26H53I	1000406-32-1	87
5			Docosane, 1-iodo-	436	C22H45I	1000406-31-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

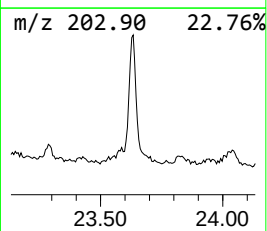
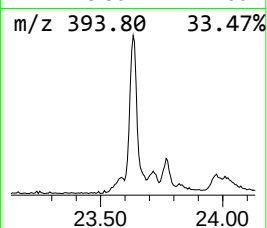
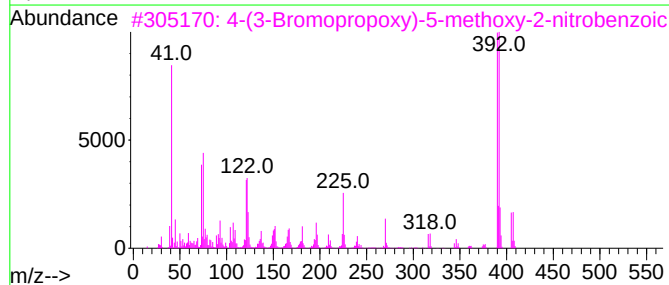
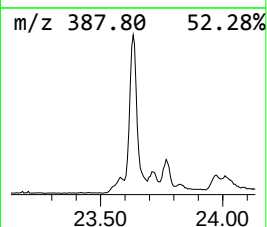
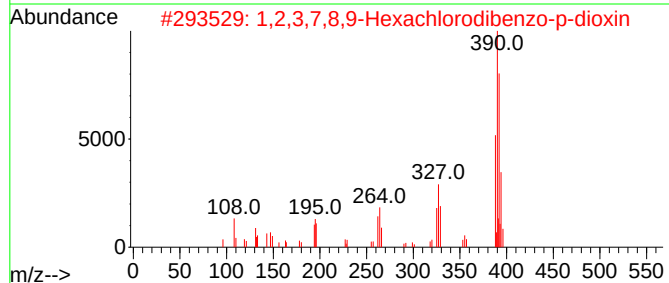
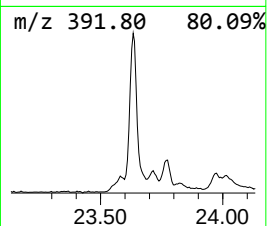
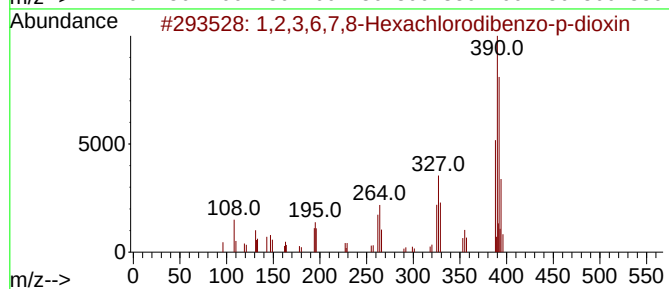
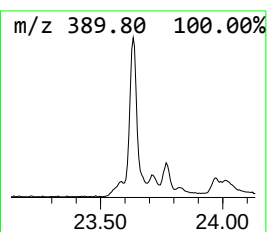
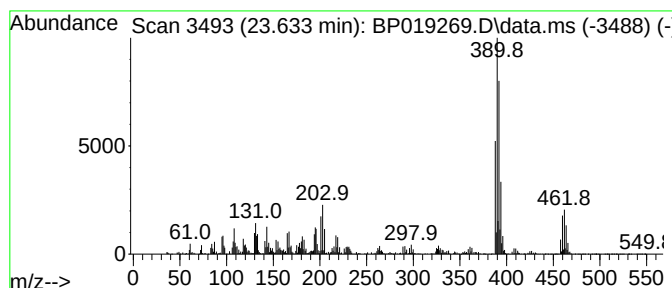
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 64 1,2,3,6,7,8-Hexachlorodiben... Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.		
23.633	20.00 ng/ul	3437340	Perylene-d12	23.692		
Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,2,3,6,7,8-Hexachlorodibenzo-p...	388	C12H2Cl6O2	057653-85-7	90	
2	1,2,3,7,8,9-Hexachlorodibenzo-p...	388	C12H2Cl6O2	019408-74-3	81	
3	4-(3-Bromopropoxy)-5-methoxy-2-n...	405	C14H20BrN06Si	1000471-26-2	38	
4	Perylene, 3-(2-naphthalenylmethyl)-	392	C31H20	086623-52-1	22	
5	N-(2,3,5,6-Tetrachlorophenyl)-2,...	425	C10H2Cl4F7NO	1000373-25-8	18	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

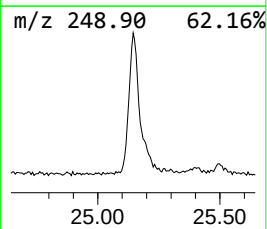
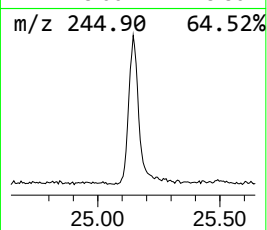
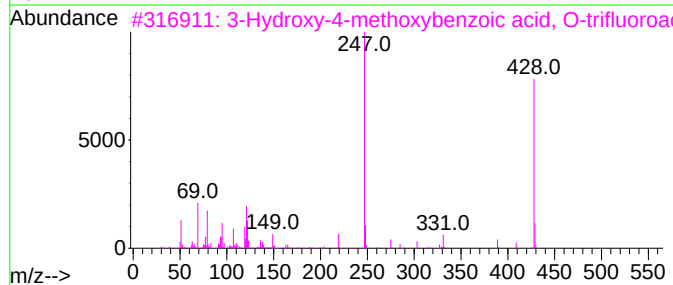
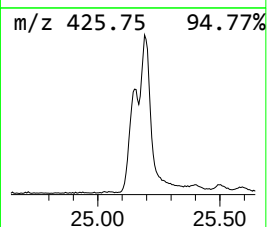
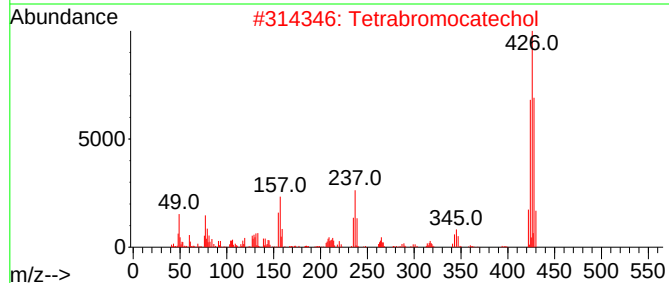
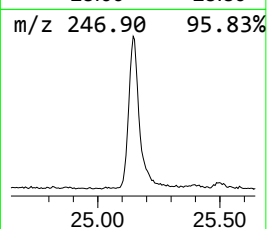
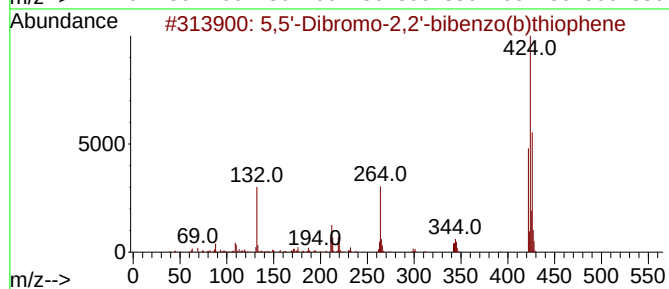
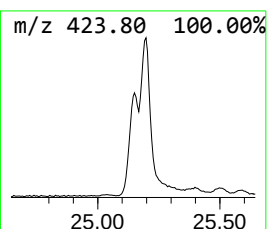
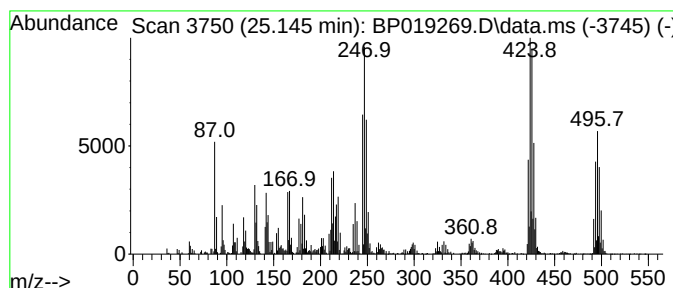
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TIC Integration Parameters: LSCINT.P

Peak Number 65 unknown-08

Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.145	18.14 ng/ul	3117210	Perylene-d12	23.692

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	5,5'-Dibromo-2,2'-bibenzo(b)thio...	422	C16H8Br2S2	007312-21-2	38		
2	Tetrabromocatechol	422	C6H2Br4O2	000488-47-1	12		
3	3-Hydroxy-4-methoxybenzoic acid,...	428	C14H9F9O5	1000447-22-1	11		
4	Acetic acid, (4-chloro-2-methylp...	424	C25H41ClO3	1000415-16-9	11		
5	2-Chloro-9,10-bis(p-methoxypheny...	424	C28H21ClO2	110904-87-5	10		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
Data File : BP019269.D
Acq On : 04 Jan 2024 18:14
Operator : MA/JU
Sample : 05951-17
Misc :
ALS Vial : 13 Sample Multiplier: 1

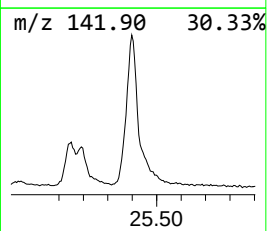
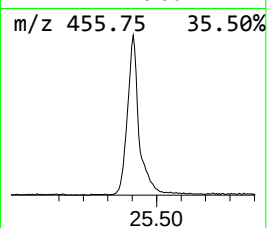
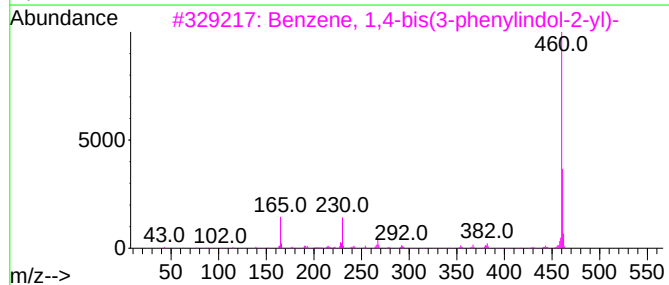
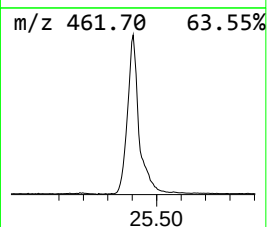
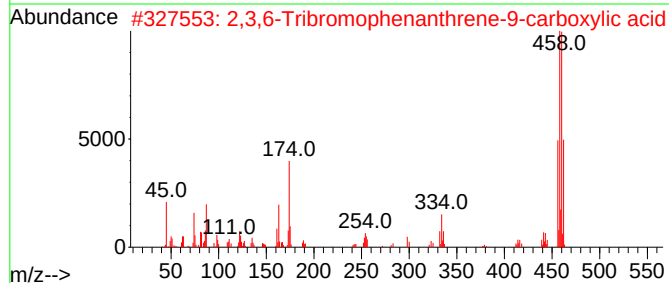
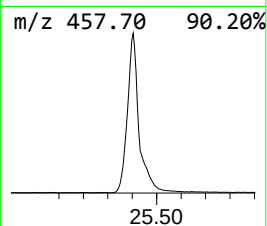
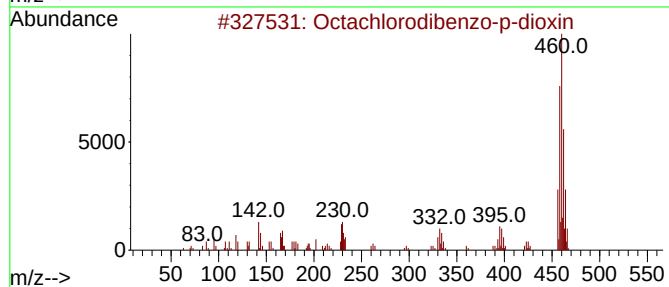
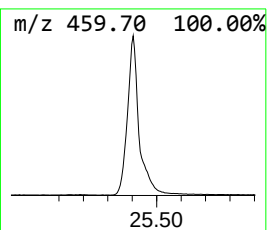
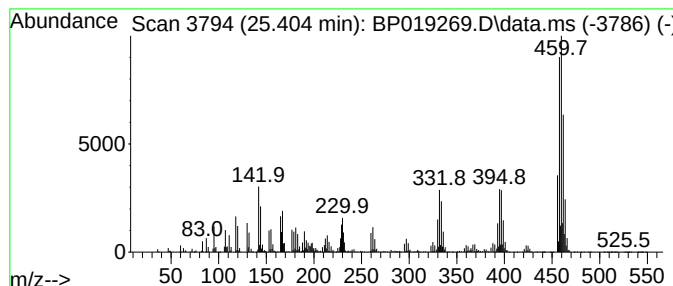
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 66 Octachlorodibenzo-p-dioxin Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.	
25.404	48.57 ng/ul	8347560	Perylene-d12	23.692	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octachlorodibenzo-p-dioxin	456	C12Cl8O2	003268-87-9	87
2	2,3,6-Tribromophenanthrene-9-car...	456	C15H7Br3O2	056988-60-4	37
3	Benzene, 1,4-bis(3-phenylindol-2...	460	C34H24N2	164936-88-3	10
4	1,3,5-Triazine-2,4,6-triamine, N...	460	C27H24N8	033933-66-3	9
5	1,1':4',1'':4'',1''':4''',1'''':...	458	C36H26	004499-83-6	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP010424\
 Data File : BP019269.D
 Acq On : 04 Jan 2024 18:14
 Operator : MA/JU
 Sample : 05951-17
 Misc :
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCF81

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP122723.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
1-Hexanol, 2-et...	7.869	59.3	ng/ul	3492050	1	7.799	1177230	20.0
Hexanoic acid, ...	9.122	7.6	ng/ul	447694	1	7.799	1177230	20.0
(DEL) Alkane: S...	10.110	5.5	ng/ul	331628	2	10.593	1196510	20.0
4-Nonanone, 2,6...	10.704	10.0	ng/ul	597880	2	10.593	1196510	20.0
1,3-Hexanediol...	10.816	13.8	ng/ul	825118	2	10.593	1196510	20.0
(DEL) Alkane: C...	10.875	5.5	ng/ul	330922	2	10.593	1196510	20.0
Neopentyl glycol	10.916	15.1	ng/ul	905845	2	10.593	1196510	20.0
(DEL) Alkane: C...	11.075	3.2	ng/ul	194126	2	10.593	1196510	20.0
(DEL) Alkane: S...	11.104	25.1	ng/ul	1498400	2	10.593	1196510	20.0
unknown-01	11.198	12.4	ng/ul	740618	2	10.593	1196510	20.0
unknown-02	11.428	12.5	ng/ul	746935	2	10.593	1196510	20.0
1H-Indene, 2,3-...	11.504	11.1	ng/ul	661229	2	10.593	1196510	20.0
(DEL) Alkane: S...	11.610	27.0	ng/ul	1614030	2	10.593	1196510	20.0
Benzene, 2-ethe...	11.698	7.4	ng/ul	441406	2	10.593	1196510	20.0
unknown-03	11.857	23.1	ng/ul	1382050	2	10.593	1196510	20.0
(DEL) Alkane: S...	12.016	114.1	ng/ul	6824520	2	10.593	1196510	20.0
unknown-04	12.381	8.4	ng/ul	505424	2	10.593	1196510	20.0
Pentadecafluoro...	12.675	51.5	ng/ul	4679580	3	14.445	1817210	20.0
(DEL) Alkane: S...	12.757	8.0	ng/ul	726039	3	14.445	1817210	20.0
(DEL) Alkane: S...	12.828	12.2	ng/ul	1107480	3	14.445	1817210	20.0
(DEL) Alkane: S...	12.975	41.8	ng/ul	3801090	3	14.445	1817210	20.0
Carbonic acid, ...	13.269	84.3	ng/ul	7654910	3	14.445	1817210	20.0
(DEL) Alkane: S...	13.357	20.8	ng/ul	1889720	3	14.445	1817210	20.0
Naphthalene, 1-...	13.475	7.6	ng/ul	691934	3	14.445	1817210	20.0
7-Methoxy-4-met...	13.498	15.5	ng/ul	1409390	3	14.445	1817210	20.0
Naphthalene, 1,...	13.610	80.5	ng/ul	7310510	3	14.445	1817210	20.0
Naphthalene, 2,...	13.769	93.0	ng/ul	8452020	3	14.445	1817210	20.0
Carbonic acid, ...	13.928	87.9	ng/ul	7987870	3	14.445	1817210	20.0
(DEL) Alkane: S...	13.963	38.4	ng/ul	3491300	3	14.445	1817210	20.0
Naphthalene, 1,...	14.004	16.2	ng/ul	1469980	3	14.445	1817210	20.0
(DEL) Alkane: S...	14.040	34.0	ng/ul	3085840	3	14.445	1817210	20.0
(DEL) Alkane: S...	14.363	346.9	ng/ul	31516600	3	14.445	1817210	20.0
Naphthalene, 1,...	14.581	20.2	ng/ul	1838980	3	14.445	1817210	20.0
Naphthalene, 2,...	14.751	42.3	ng/ul	3838470	3	14.445	1817210	20.0
Naphthalene, 1,...	14.857	136.2	ng/ul	12372800	3	14.445	1817210	20.0
unknown-05	14.934	125.5	ng/ul	11398600	3	14.445	1817210	20.0
(DEL) Alkane: S...	14.975	80.6	ng/ul	7322830	3	14.445	1817210	20.0
2-Bromotetradecane	15.045	29.4	ng/ul	2674160	3	14.445	1817210	20.0
unknown-06	15.286	85.5	ng/ul	7765700	3	14.445	1817210	20.0
(DEL) Alkane: S...	15.328	177.8	ng/ul	16157900	3	14.445	1817210	20.0
unknown-07	15.728	171.7	ng/ul	15602900	3	14.445	1817210	20.0
(DEL) Alkane: S...	15.881	7.3	ng/ul	10776500	4	17.275	29435600	20.0
(DEL) Alkane: S...	15.939	6.3	ng/ul	9234510	4	17.275	29435600	20.0
(DEL) Alkane: S...	16.204	20.8	ng/ul	30566700	4	17.275	29435600	20.0
Phenanthrene, 2...	19.027	9.2	ng/ul	13587900	4	17.275	29435600	20.0
(DEL) Alkane: S...	20.574	28.3	ng/ul	13331900	5	21.333	9425080	20.0
(DEL) Alkane: S...	21.027	13.6	ng/ul	6428840	5	21.333	9425080	20.0
(DEL) Alkane: S...	21.457	14.8	ng/ul	6953890	5	21.333	9425080	20.0
1,2,3,6,7,8-Hex...	23.633	20.0	ng/ul	3437340	6	23.692	3437340	20.0
unknown-08	25.145	18.1	ng/ul	3117210	6	23.692	3437340	20.0
Octachlorodiben...	25.404	48.6	ng/ul	8347560	6	23.692	3437340	20.0