

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
 Data File : BP022160.D
 Acq On : 02 Oct 2024 13:25
 Operator : RC/JU
 Sample : P4016-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCVW3

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-BP092424.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022160.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	5.752	464	470	479	rVB2	1979495	3022485	1.72%	0.479%
2	6.711	622	633	638	rBV3	1617987	3760301	2.14%	0.596%
3	6.769	638	643	646	rBV	1741877	2521219	1.44%	0.399%
4	6.811	646	650	654	rVB	2855433	4094819	2.33%	0.649%
5	6.869	654	660	667	rBV4	2913876	5692563	3.24%	0.902%
6	6.940	667	672	677	rVB	1654874	2308100	1.31%	0.366%
7	7.099	694	699	703	rBV	1761051	2908566	1.66%	0.461%
8	7.140	703	706	713	rVB	1515832	2139983	1.22%	0.339%
9	7.346	734	741	749	rBV2	9422700	19275748	10.98%	3.053%
10	7.622	777	788	793	rVV5	805045	2668677	1.52%	0.423%
11	7.687	793	799	805	rVV2	2437898	4307404	2.45%	0.682%
12	7.769	805	813	817	rVV3	1269539	3032514	1.73%	0.480%
13	7.816	817	821	829	rVV3	1857902	3627030	2.07%	0.575%
14	7.922	836	839	842	rVV	1272559	1898900	1.08%	0.301%
15	7.987	847	850	857	rVV3	1072220	2094291	1.19%	0.332%
16	8.128	869	874	878	rBV	1864715	2979603	1.70%	0.472%
17	8.181	878	883	887	rVV2	1450871	3084585	1.76%	0.489%
18	8.234	887	892	898	rVV2	2450112	5374464	3.06%	0.851%
19	8.287	898	901	905	rVV	1690163	2311998	1.32%	0.366%
20	8.334	905	909	926	rVB4	3200747	9824434	5.60%	1.556%
21	8.463	926	931	934	rBV	2174815	3627612	2.07%	0.575%
22	8.493	934	936	945	rVB2	1841318	2964699	1.69%	0.470%
23	8.646	957	962	965	rBV	1640724	2513771	1.43%	0.398%
24	8.693	965	970	977	rVB2	2148689	4033713	2.30%	0.639%
25	8.793	977	987	993	rBV3	2616064	5834640	3.32%	0.924%
26	8.922	998	1009	1018	rVB	9300760	17211002	9.80%	2.726%
27	9.040	1018	1029	1035	rBV7	1967011	5825672	3.32%	0.923%
28	9.116	1035	1042	1048	rBV3	1319568	3205746	1.83%	0.508%
29	9.469	1095	1102	1109	rVB4	968460	2163437	1.23%	0.343%
30	9.558	1109	1117	1122	rBV5	1089564	2678259	1.53%	0.424%
31	9.616	1122	1127	1131	rBV3	1347885	2613433	1.49%	0.414%
32	9.758	1148	1151	1156	rVB	1715608	2410140	1.37%	0.382%
33	9.822	1156	1162	1166	rBV3	1237074	2011476	1.15%	0.319%
34	10.005	1187	1193	1198	rBV2	836880	1829208	1.04%	0.290%
35	10.163	1215	1220	1226	rBV2	1330147	2242662	1.28%	0.355%
36	10.475	1260	1273	1278	rBV3	11319433	28665248	16.33%	4.541%

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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-BP092424.MA.M
 Title : SVOA CALIBRATION

37	10.587	1286	1292	1297	rBV2	5362922	8938072	5.09%	1.416%
38	10.846	1330	1336	1344	rBV	4504731	7767904	4.43%	1.230%
39	10.975	1349	1358	1364	rVV	3720501	6954053	3.96%	1.102%
40	11.493	1440	1446	1452	rBV	3494902	6035351	3.44%	0.956%

41	11.563	1453	1458	1465	rVB4	1274451	2541353	1.45%	0.403%
42	11.728	1478	1486	1494	rVB2	2613927	4611995	2.63%	0.731%
43	11.893	1507	1514	1517	rBV	3857111	6338894	3.61%	1.004%
44	11.928	1517	1520	1525	rVB	2448313	3172235	1.81%	0.502%
45	12.134	1548	1555	1563	rVB2	4108040	7789484	4.44%	1.234%

46	12.304	1578	1584	1588	rBV3	4247011	7363919	4.20%	1.166%
47	12.563	1613	1628	1636	rBV5	1294907	4463228	2.54%	0.707%
48	12.646	1637	1642	1648	rVV6	942700	2067223	1.18%	0.327%
49	12.710	1648	1653	1662	rVV5	804366	2125922	1.21%	0.337%
50	12.887	1679	1683	1688	rVB2	1266987	2063458	1.18%	0.327%

51	13.140	1721	1726	1732	rBV	4297125	6666303	3.80%	1.056%
52	13.369	1758	1765	1771	rVB4	2634070	5798206	3.30%	0.918%
53	13.498	1777	1787	1792	rBV4	2737660	6573841	3.75%	1.041%
54	13.646	1806	1812	1815	rBV2	2262488	4349967	2.48%	0.689%
55	13.804	1831	1839	1843	rBV4	1386250	3092450	1.76%	0.490%

56	13.869	1843	1850	1854	rVB6	1340073	2601783	1.48%	0.412%
57	13.963	1862	1866	1868	rBV3	1392705	2252938	1.28%	0.357%
58	13.981	1868	1869	1877	rVB3	1798691	2596820	1.48%	0.411%
59	14.051	1877	1881	1885	rBV2	2835570	3639442	2.07%	0.576%
60	14.128	1891	1894	1907	rVB9	869103	2453652	1.40%	0.389%

61	14.240	1907	1913	1917	rBV	6028255	8548470	4.87%	1.354%
62	14.287	1917	1921	1924	rBV4	1475308	2714975	1.55%	0.430%
63	14.322	1924	1927	1932	rVV	1363411	2425729	1.38%	0.384%
64	14.398	1935	1940	1947	rVV6	2171575	4055528	2.31%	0.642%
65	14.463	1947	1951	1955	rVB2	1927403	2654397	1.51%	0.420%

66	14.540	1960	1964	1968	rBV4	1025779	1814808	1.03%	0.287%
67	14.734	1994	1997	2001	rVB	2129161	2782519	1.59%	0.441%
68	14.798	2002	2008	2013	rVB2	1587701	2572514	1.47%	0.407%
69	14.857	2014	2018	2023	rBV4	1179955	2746071	1.56%	0.435%
70	14.981	2030	2039	2044	rVB2	13507738	25657671	14.62%	4.064%

71	15.098	2055	2059	2062	rBV2	2671094	3677836	2.10%	0.583%
72	15.193	2067	2075	2080	rVB3	4508750	7821211	4.46%	1.239%
73	15.445	2113	2118	2121	rVV5	2286141	3929132	2.24%	0.622%
74	15.487	2121	2125	2133	rVB2	2216646	3672879	2.09%	0.582%
75	15.616	2142	2147	2156	rVB5	1956341	4933192	2.81%	0.781%

76	15.775	2167	2174	2177	rVV3	903457	1927194	1.10%	0.305%
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Integration Parameters: LSCINT.P

Integrator: RTE

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Stop Thrs : 0

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 Title : SVOA CALIBRATION

77	15.975	2202	2208	2215	rBV5	1075018	2812917	1.60%	0.446%
78	16.075	2220	2225	2233	rBV2	4768513	8814657	5.02%	1.396%
79	16.834	2337	2354	2357	rBV3	40056925	175535879	100.00%	27.805%
80	16.887	2360	2363	2367	rVB	4268583	4865604	2.77%	0.771%
81	16.939	2368	2372	2376	rBV	1620916	2354638	1.34%	0.373%
82	17.145	2401	2407	2410	rBV	1787545	2856890	1.63%	0.453%
83	17.187	2410	2414	2418	rVV	1034951	2114209	1.20%	0.335%
84	17.234	2418	2422	2425	rVV	2096654	2845753	1.62%	0.451%
85	17.445	2451	2458	2462	rBV5	1396300	2751797	1.57%	0.436%
86	17.645	2487	2492	2497	rVB	3810655	5458603	3.11%	0.865%
87	17.828	2519	2523	2527	rVB	5167811	6029761	3.44%	0.955%
88	18.063	2560	2563	2569	rVB2	1728476	2512086	1.43%	0.398%
89	18.369	2610	2615	2620	rVB	2757824	4004936	2.28%	0.634%
90	19.039	2721	2729	2735	rVV3	3560367	8127659	4.63%	1.287%
91	19.181	2749	2753	2758	rBV2	2086407	2663369	1.52%	0.422%
92	19.598	2820	2824	2827	rBV	1472969	2397249	1.37%	0.380%
93	19.628	2827	2829	2839	rVB2	1957177	2614322	1.49%	0.414%
94	20.180	2920	2923	2929	rVB2	1862543	2462357	1.40%	0.390%
95	21.227	3097	3101	3115	rVB2	2585059	4277519	2.44%	0.678%
96	21.551	3151	3156	3164	rVB2	1495739	2478771	1.41%	0.393%
97	21.792	3193	3197	3202	rVB	1253848	1973999	1.12%	0.313%
98	22.416	3297	3303	3312	rVB2	2275023	4548964	2.59%	0.721%
99	24.616	3670	3677	3689	rVB	1029127	2601379	1.48%	0.412%
100	24.845	3709	3716	3726	rVB2	1259649	3258133	1.86%	0.516%

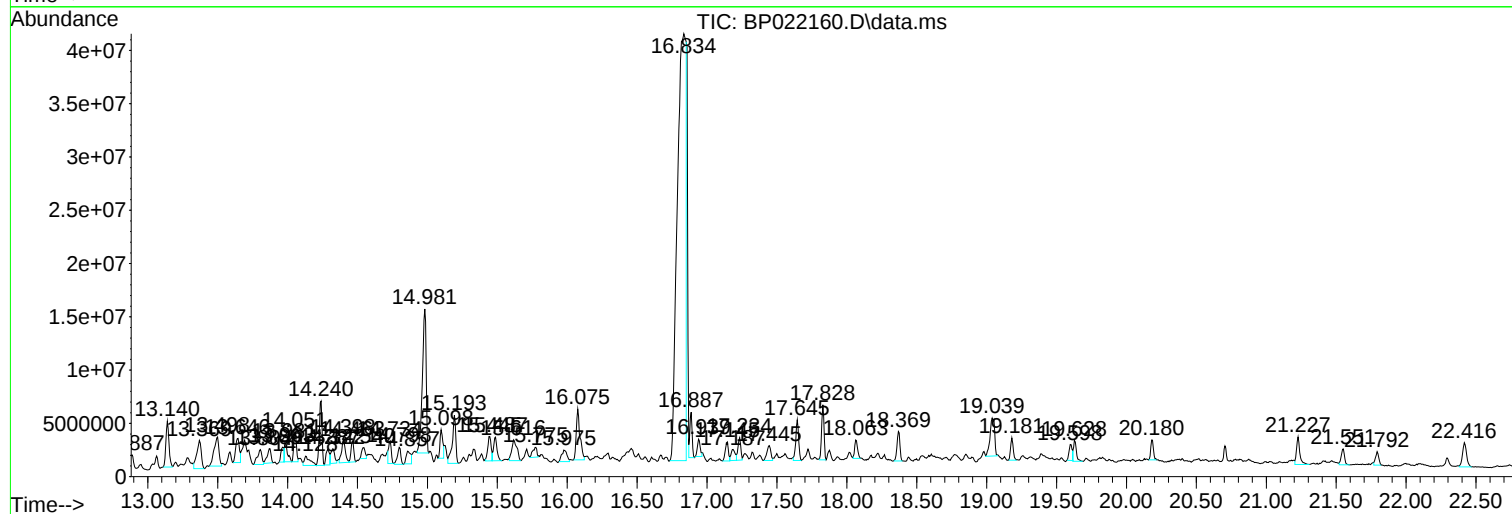
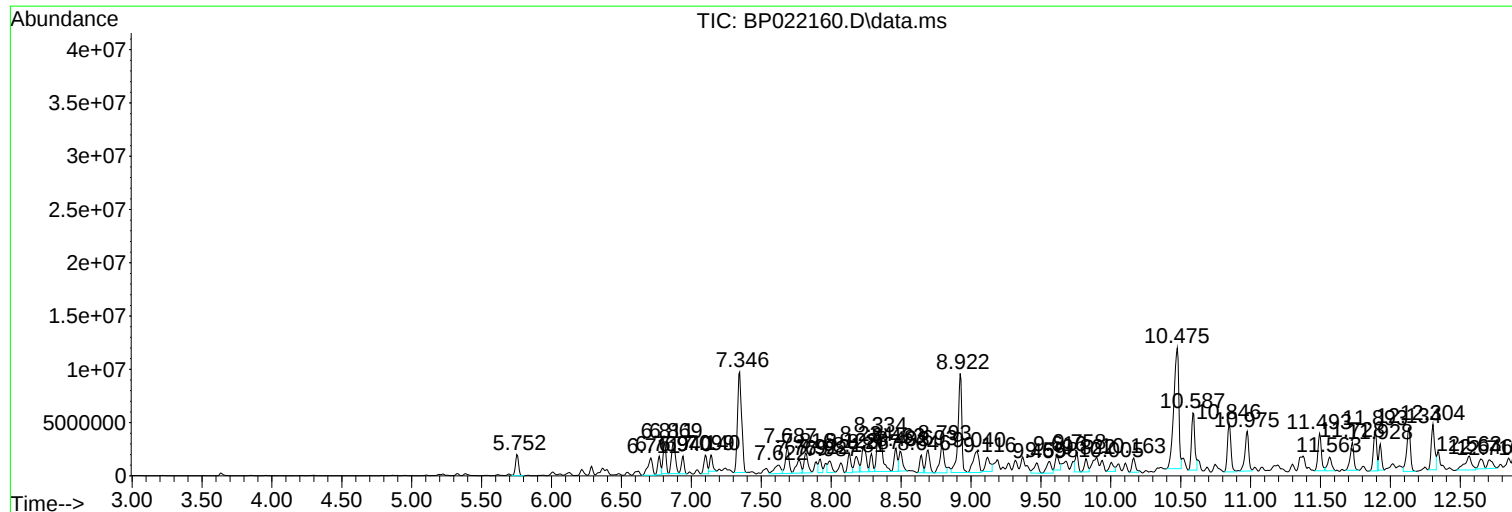
Sum of corrected areas: 631312472

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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



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ALS Vial : 5 Sample Multiplier: 1

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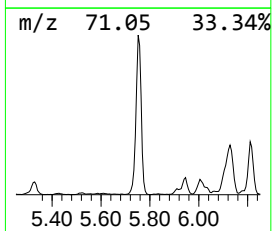
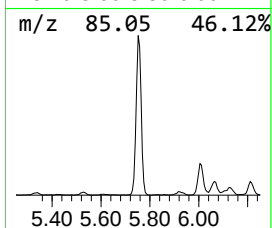
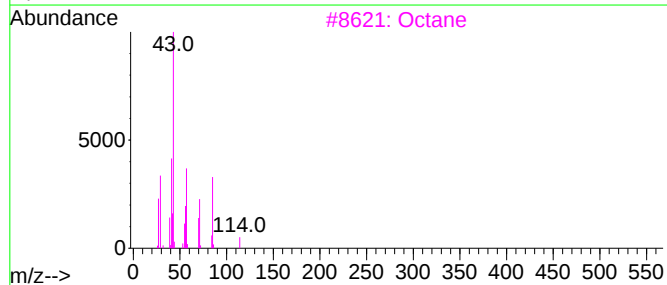
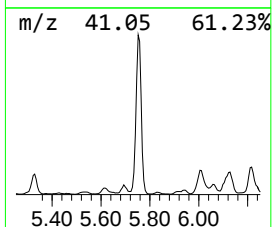
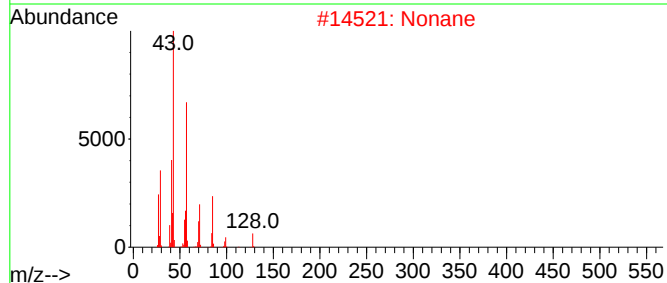
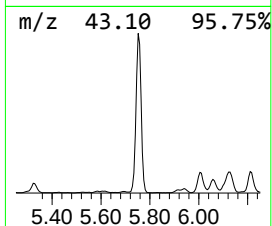
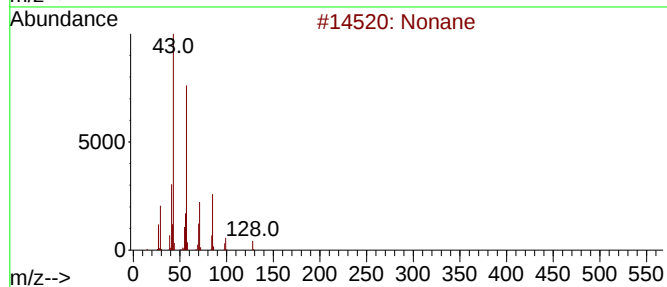
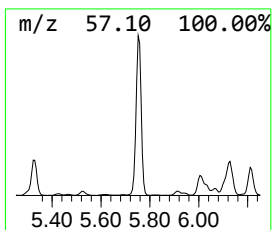
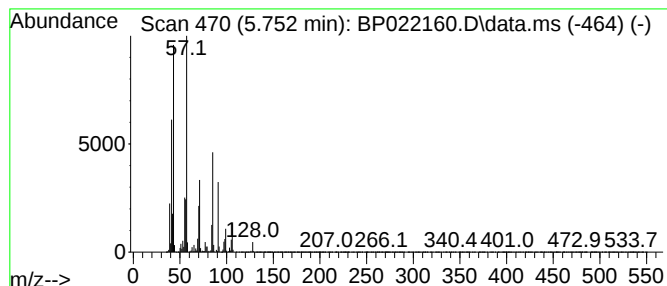
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 45

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.752	14.03 ng/ul	3022490	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane	128	C9H20	000111-84-2	93
2			Nonane	128	C9H20	000111-84-2	93
3			Octane	114	C8H18	000111-65-9	64
4			Octane	114	C8H18	000111-65-9	59
5			Sulfurous acid, hexyl heptyl ester	264	C13H28O3S	1000309-12-9	50



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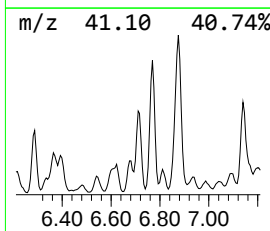
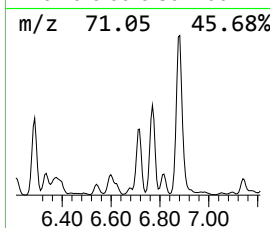
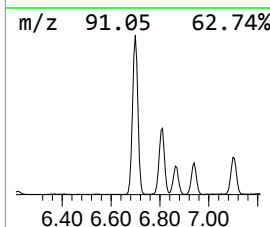
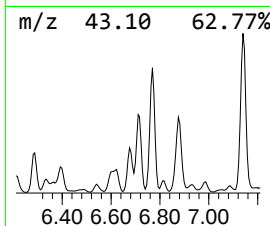
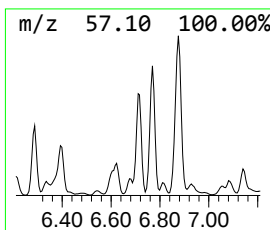
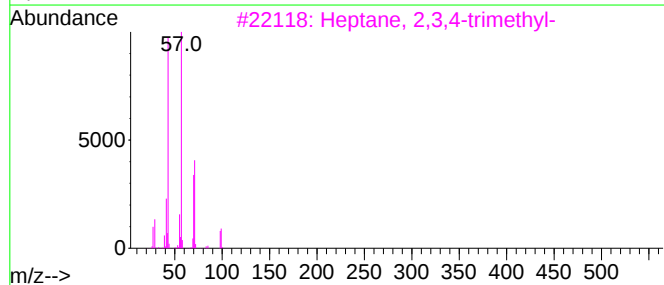
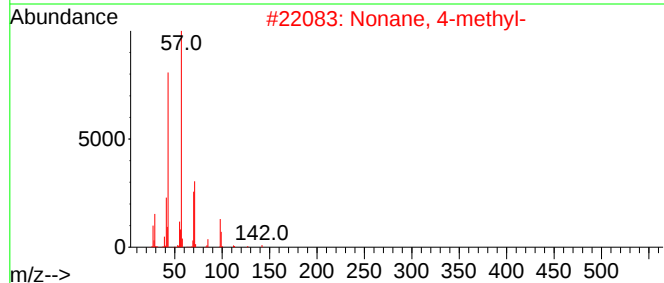
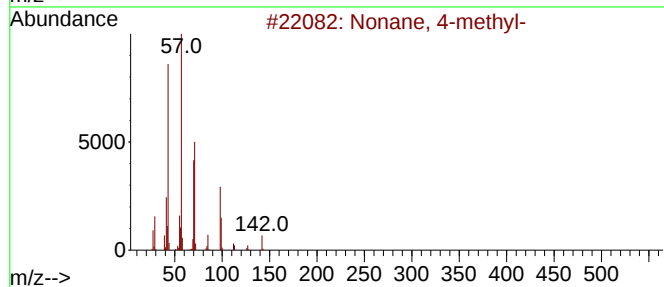
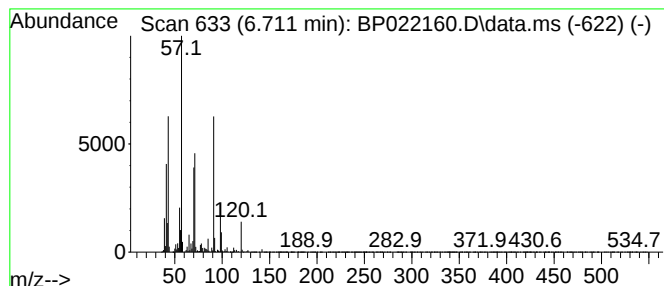
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 36

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.711	17.46 ng/ul	3760300	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 4-methyl-	142	C10H22	017301-94-9	72
2			Nonane, 4-methyl-	142	C10H22	017301-94-9	70
3			Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	59
4			Nonane, 4-methyl-	142	C10H22	017301-94-9	53
5			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53



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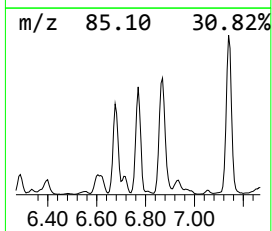
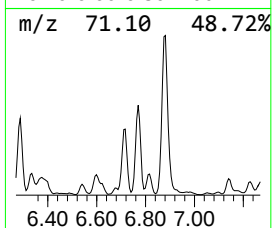
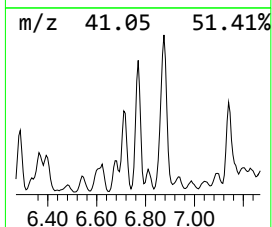
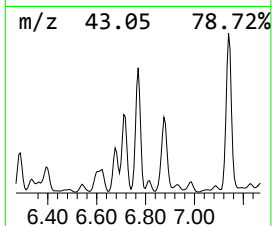
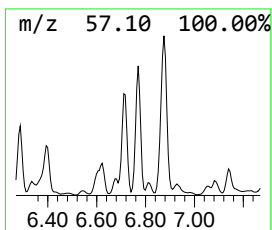
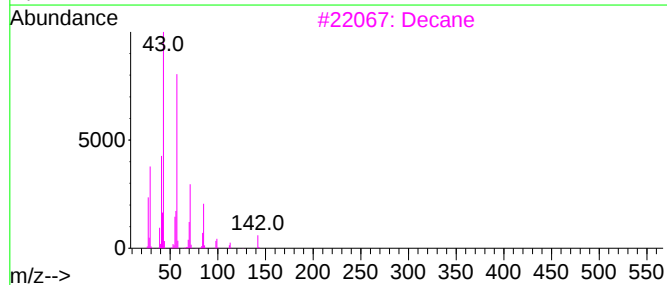
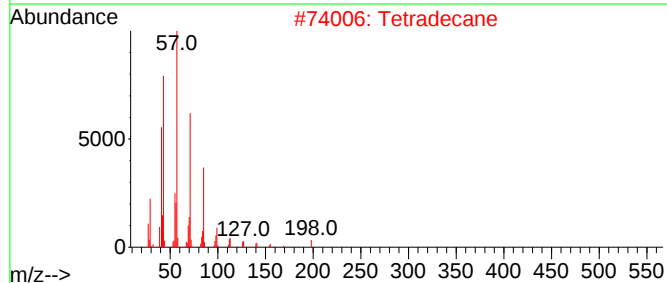
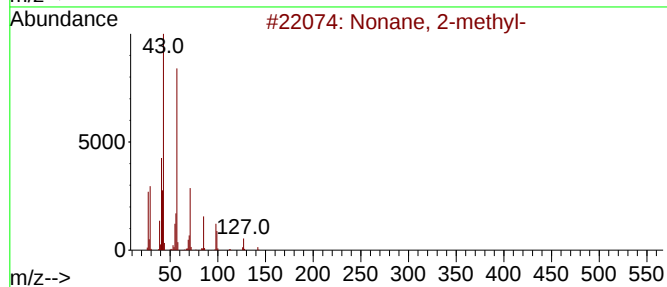
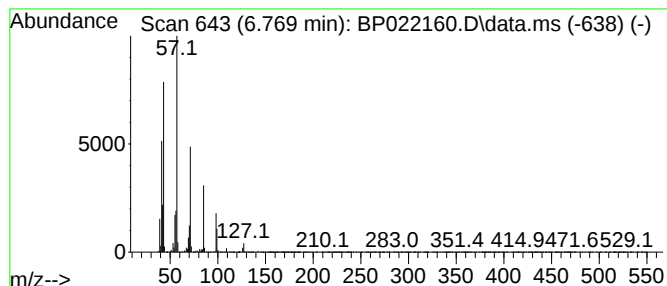
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 51

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.769	11.71 ng/ul	2521220	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonane, 2-methyl-	142	C10H22	000871-83-0	94
2			Tetradecane	198	C14H30	000629-59-4	72
3			Decane	142	C10H22	000124-18-5	64
4			Tridecane	184	C13H28	000629-50-5	59
5			Nonane, 2-methyl-	142	C10H22	000871-83-0	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

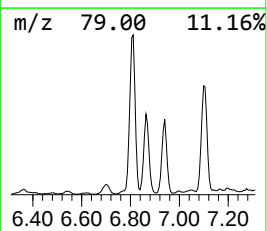
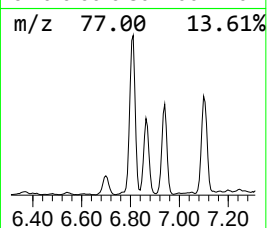
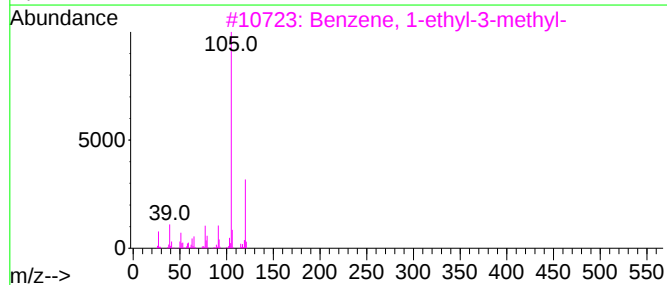
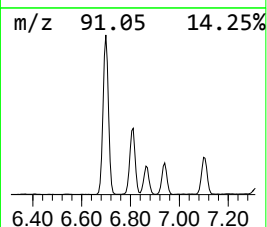
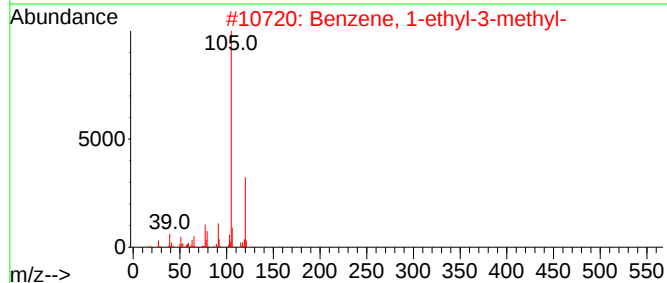
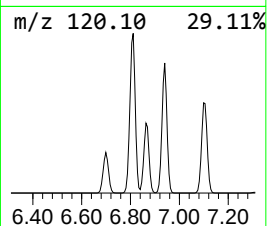
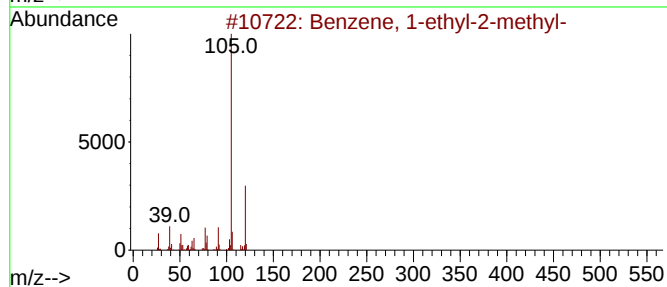
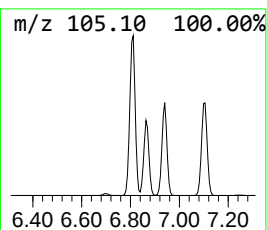
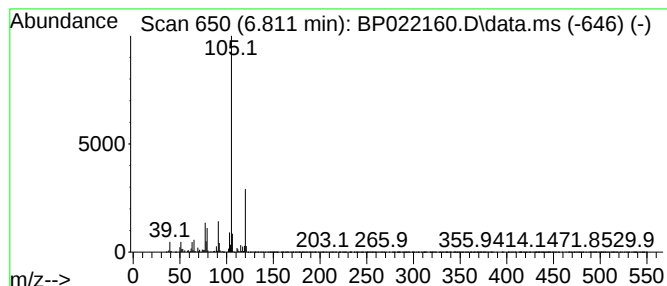
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 32

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.811	19.01 ng/ul	4094820	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	95
2			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
3			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	94
4			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5			Mesitylene	120	C9H12	000108-67-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

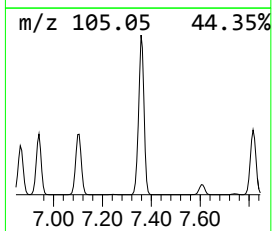
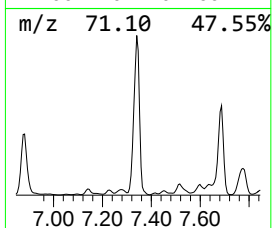
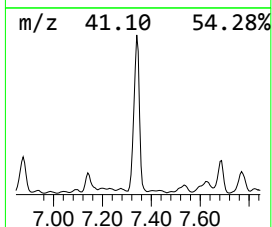
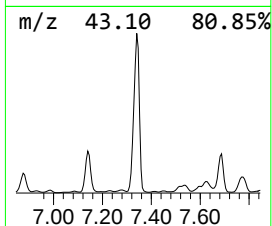
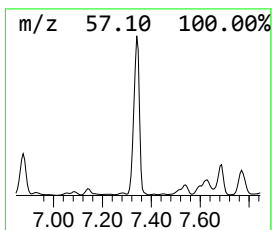
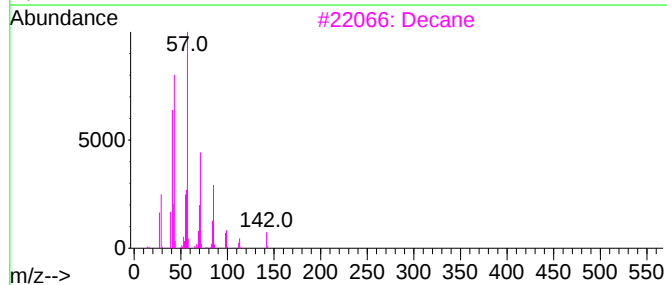
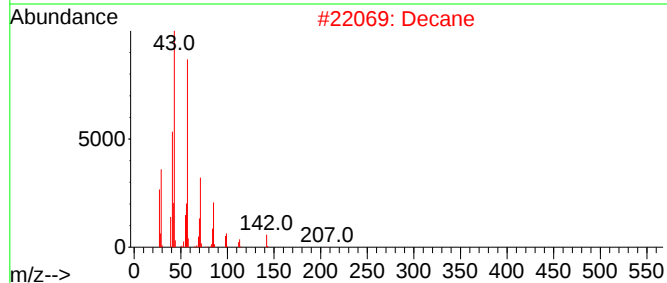
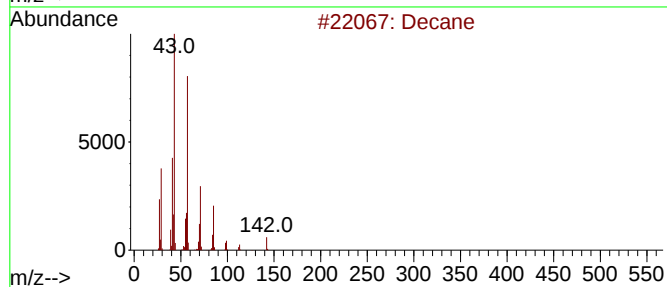
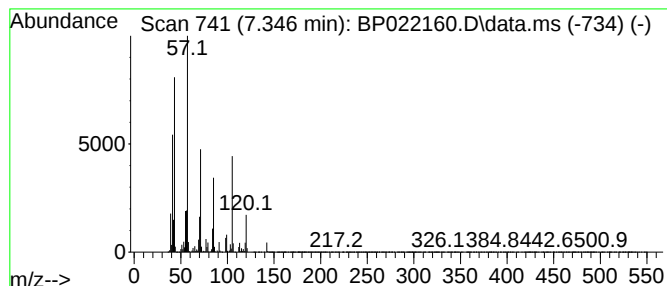
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.346	89.50 ng/ul	19275700	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane			142	C10H22	000124-18-5	93
2	Decane			142	C10H22	000124-18-5	87
3	Decane			142	C10H22	000124-18-5	87
4	Decane			142	C10H22	000124-18-5	87
5	Undecane			156	C11H24	001120-21-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

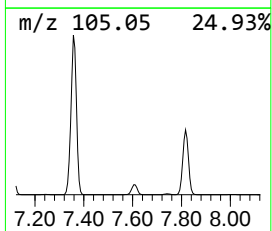
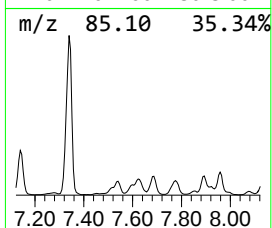
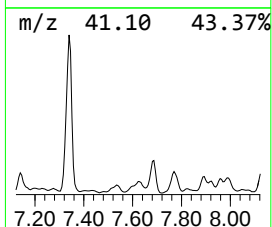
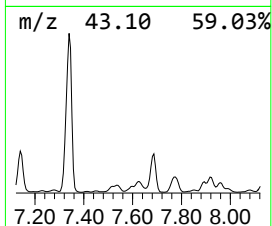
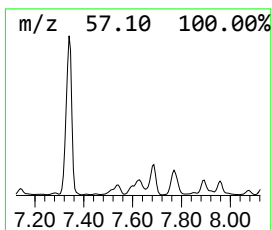
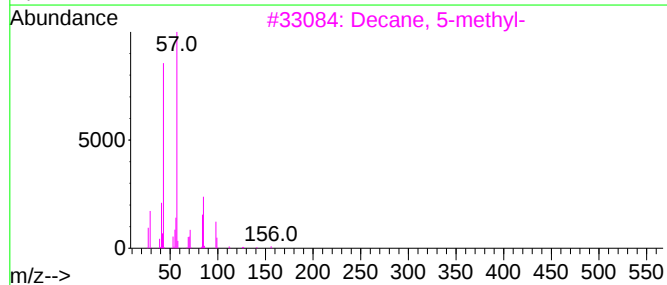
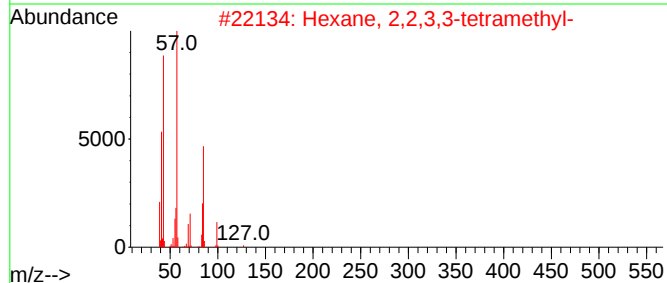
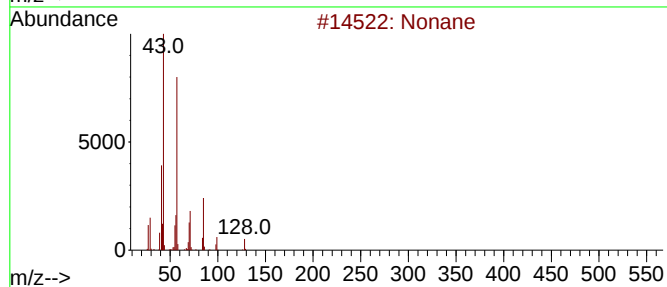
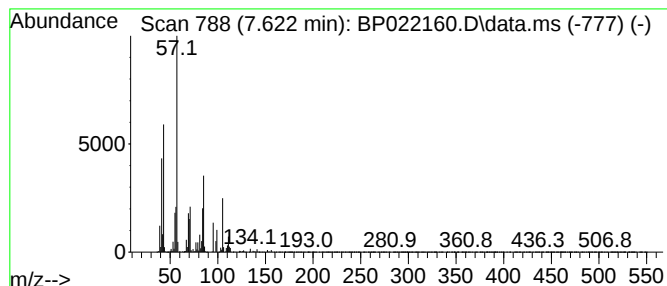
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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.	
7.622	12.39 ng/ul	2668680	1,4-Dichlorobenzene-d4	7.699	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane	128	C9H20	000111-84-2	64
2	Hexane, 2,2,3,3-tetramethyl-	142	C10H22	013475-81-5	53
3	Decane, 5-methyl-	156	C11H24	013151-35-4	47
4	Decane, 2,5,6-trimethyl-	184	C13H28	062108-23-0	47
5	Decane	142	C10H22	000124-18-5	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Sample : P4016-12
Misc :
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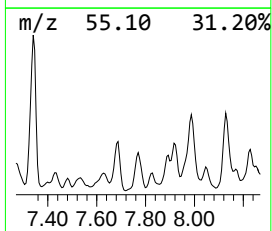
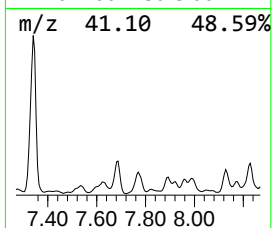
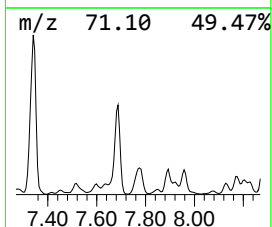
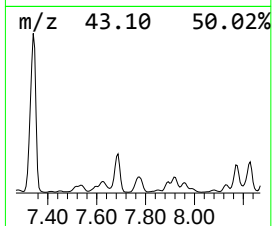
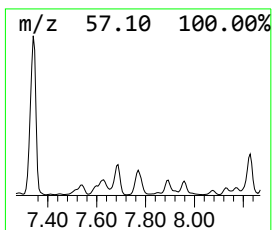
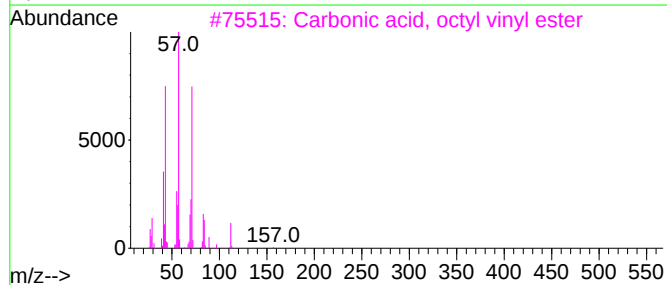
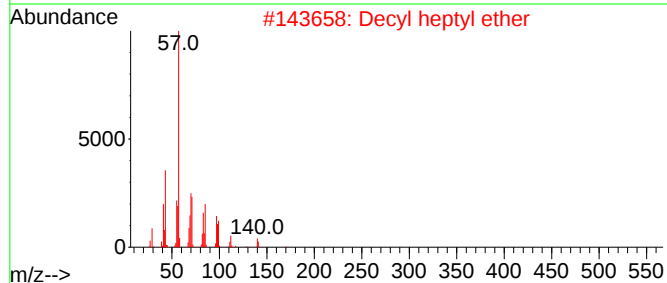
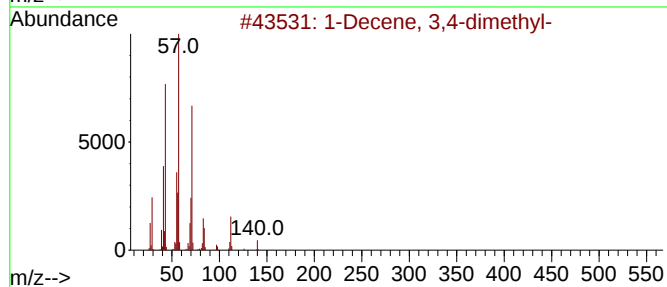
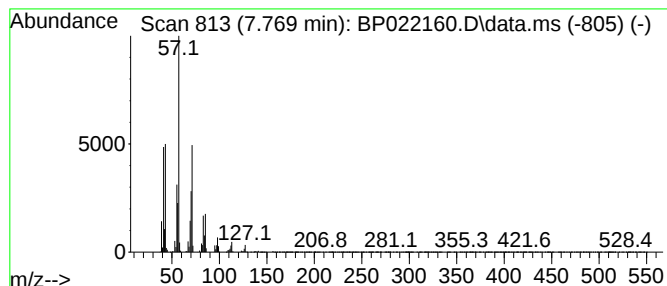
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 44

R.T.	EstConc	Area	Relative to ISTD	R.T.	
7.769	14.08 ng/ul	3032510	1,4-Dichlorobenzene-d4	7.699	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decene, 3,4-dimethyl-	168	C12H24	050871-03-9	64
2	Decyl heptyl ether	256	C17H36O	1000406-39-1	64
3	Carbonic acid, octyl vinyl ester	200	C11H20O3	1000383-25-5	53
4	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	53
5	Heptadecane, 2,6,10,14-tetramethyl-	296	C21H44	018344-37-1	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
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Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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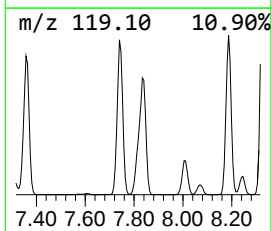
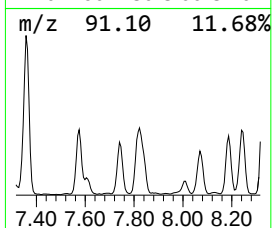
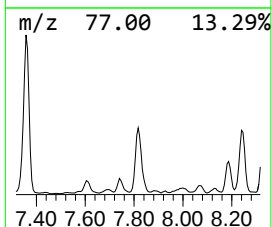
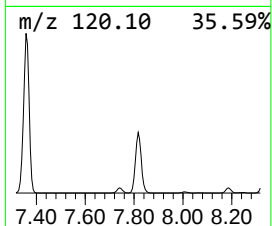
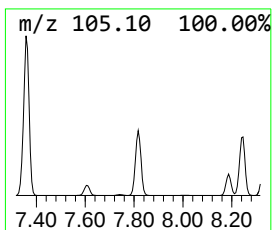
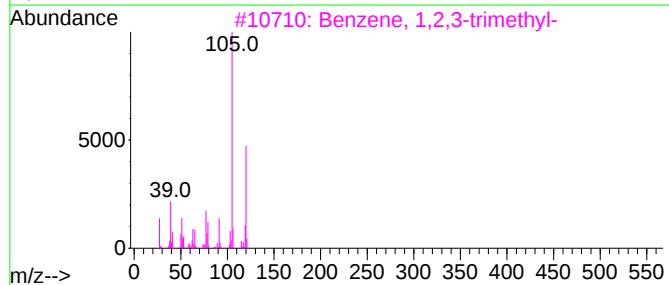
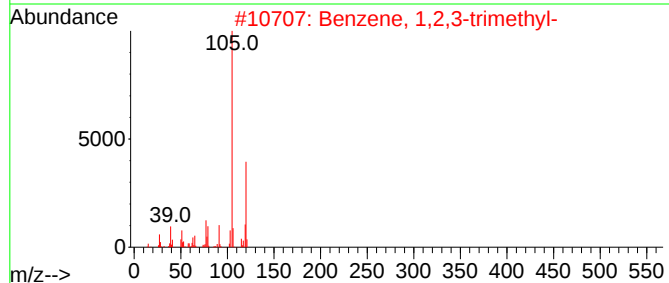
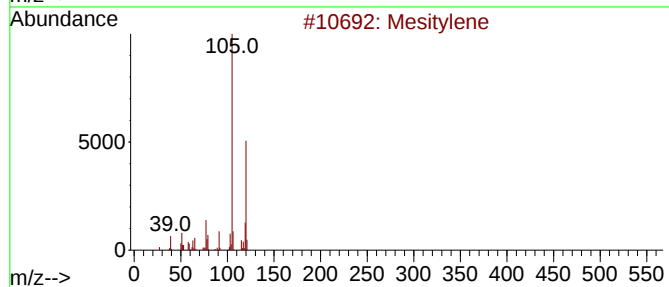
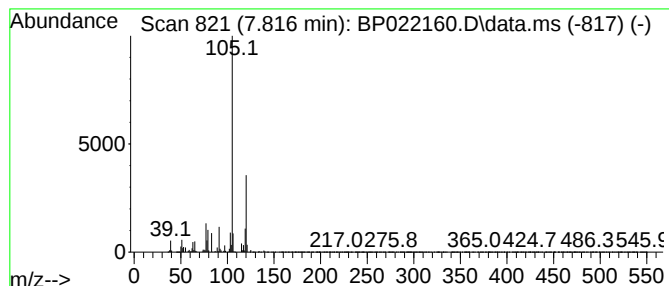
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Mesitylene Concentration Rank 40

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.816	16.84 ng/ul	3627030	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	95
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
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Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

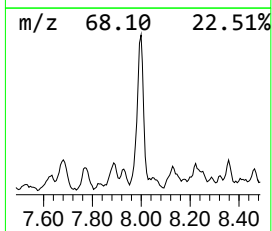
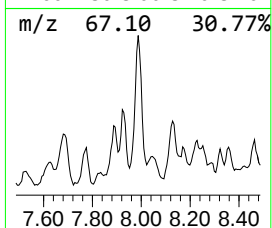
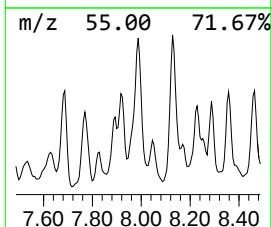
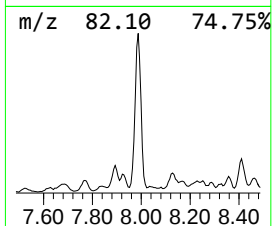
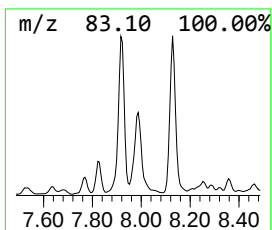
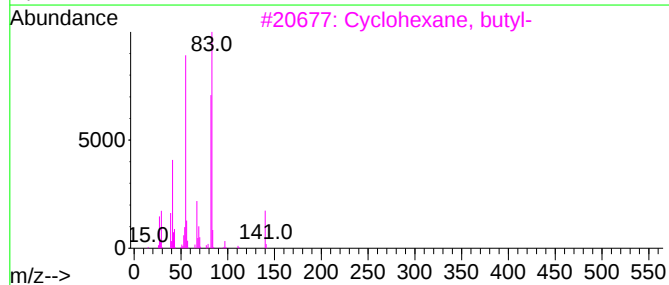
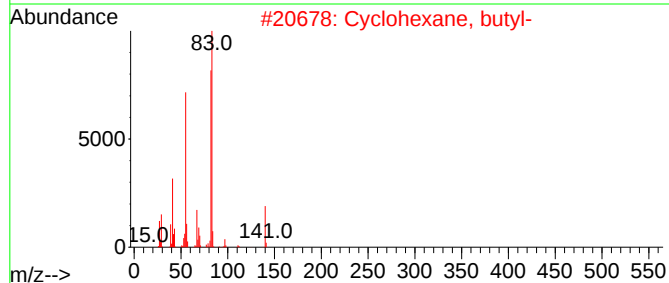
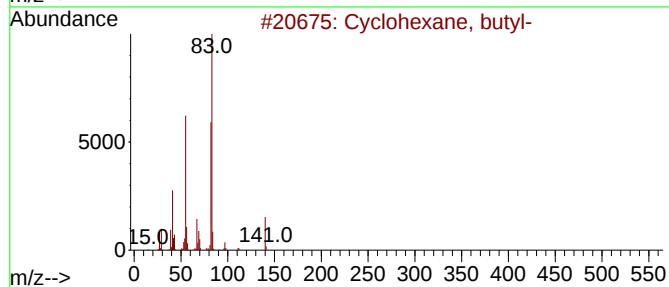
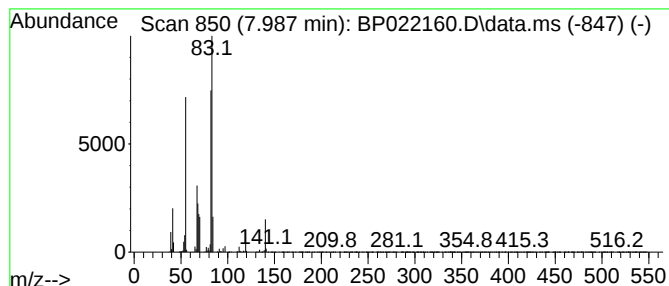
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Cyclic7.987 Concentration Rank 55

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	9.72 ng/ul	2094290	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexane, butyl-	140	C10H20	001678-93-9	80
2			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
3			Cyclohexane, butyl-	140	C10H20	001678-93-9	72
4			Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	64
5			Cyclohexane, propyl-	126	C9H18	001678-92-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

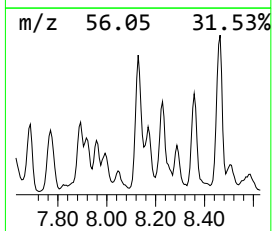
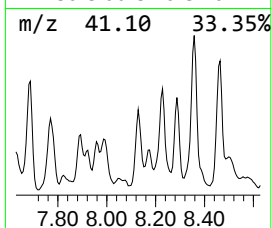
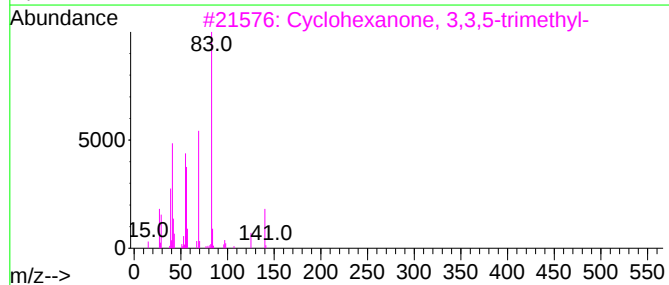
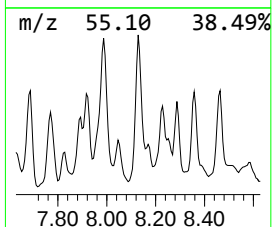
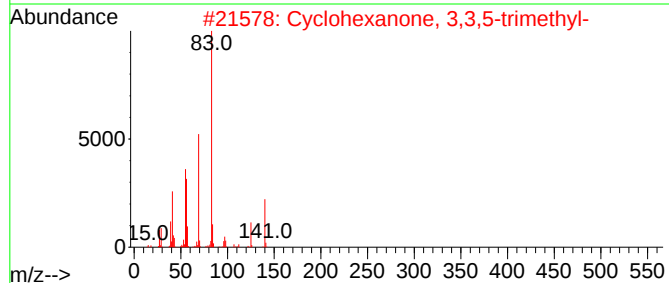
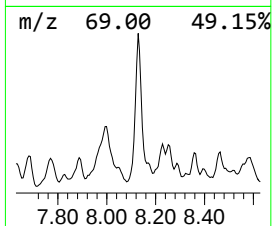
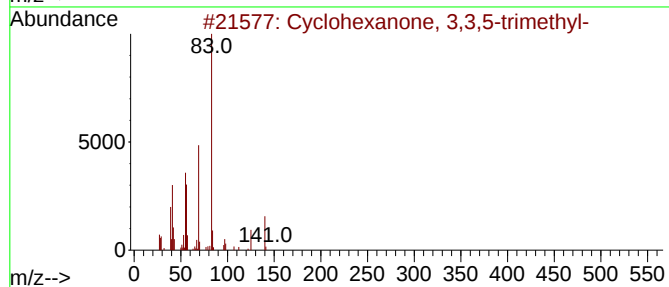
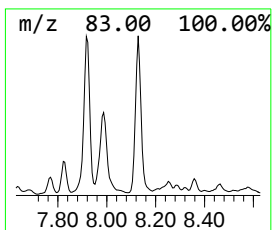
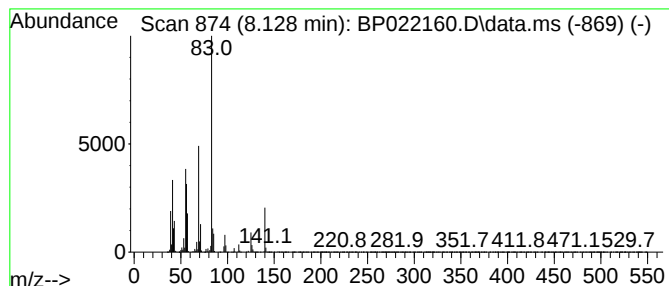
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Cyclohexanone, 3,3,5-trimet... Concentration Rank 46

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.128	13.83 ng/ul	2979600	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	94
2			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	93
3			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	87
4			Cyclohexanone, 3,3,5-trimethyl-	140	C9H16O	000873-94-9	83
5			Cyclohexane, (2-methylpropyl)-	140	C10H20	001678-98-4	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

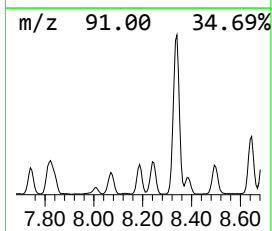
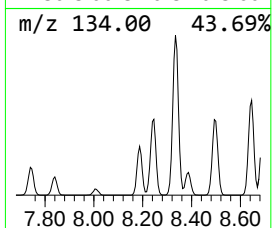
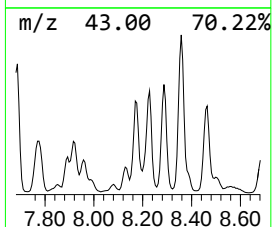
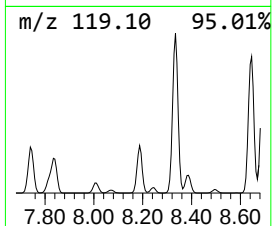
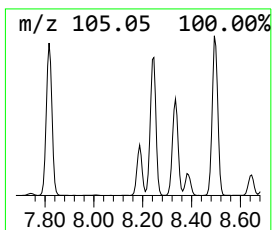
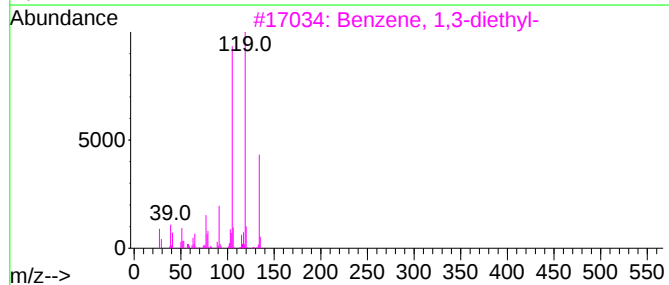
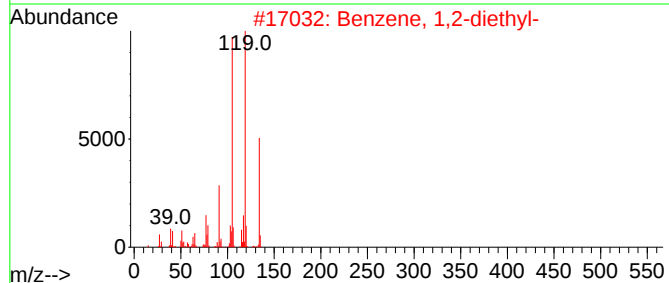
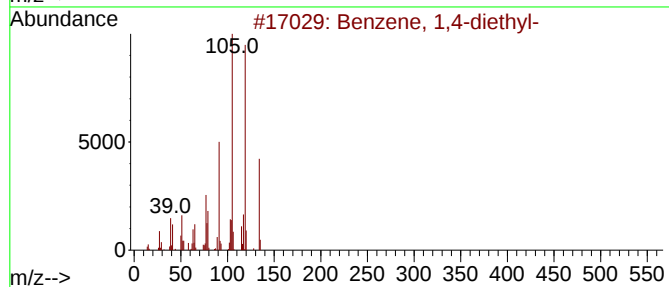
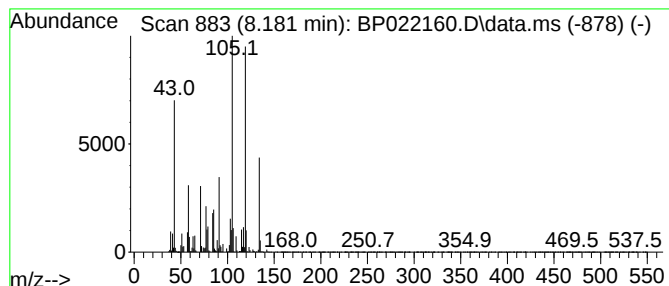
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzene, 1,4-diethyl- Concentration Rank 43

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.181	14.32 ng/ul	3084590	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	94
4			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	94
5			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

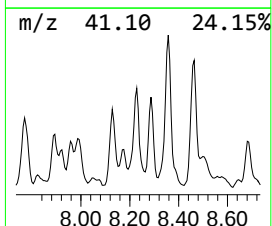
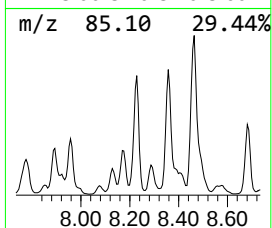
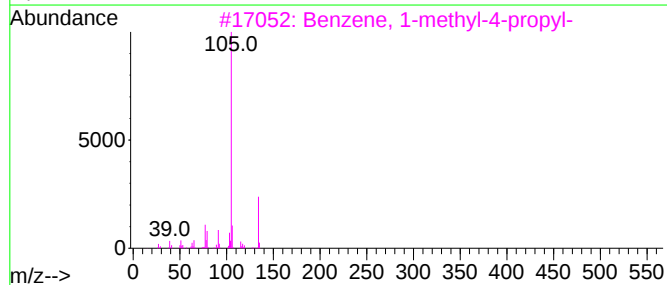
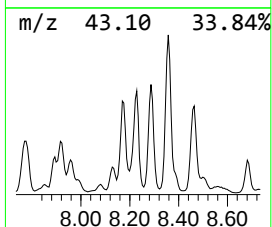
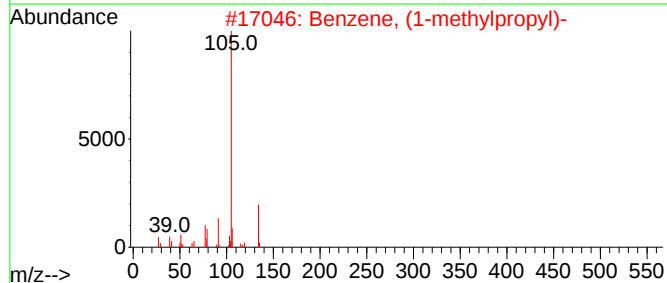
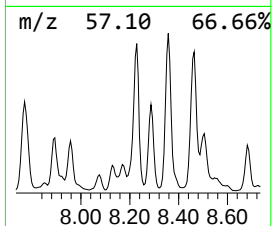
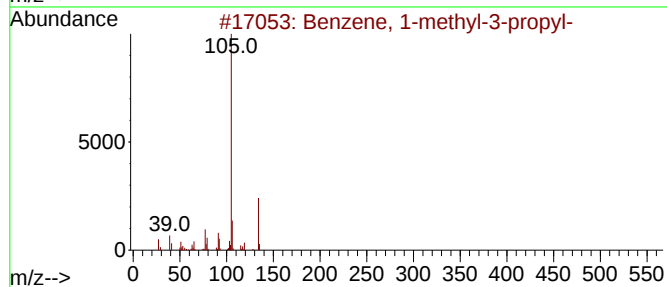
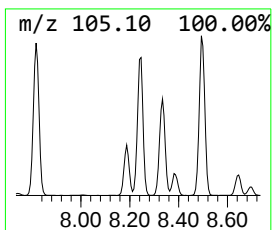
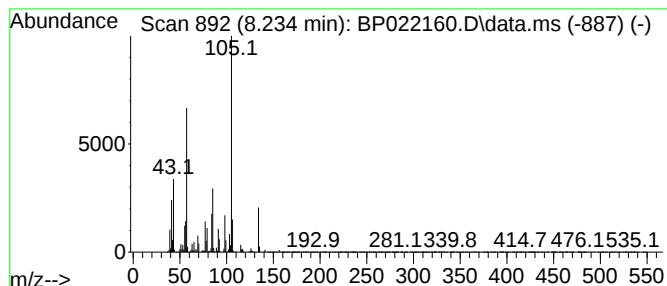
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzene, 1-methyl-3-propyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.234	24.95 ng/ul	5374460	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	55
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	50
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	50
4			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	50
5			Benzene, 1-methyl-3-propyl-	134	C10H14	001074-43-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

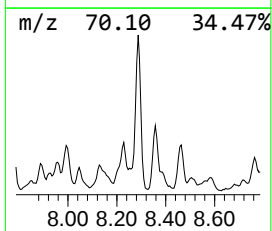
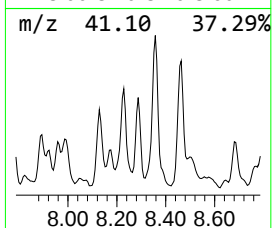
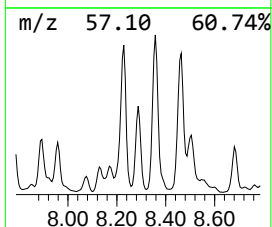
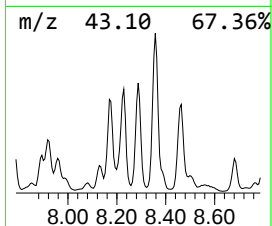
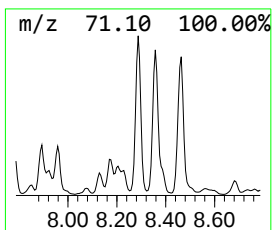
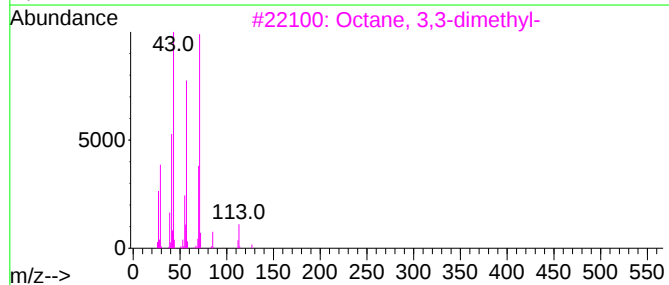
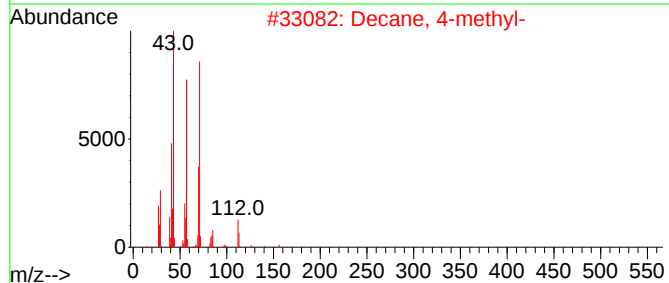
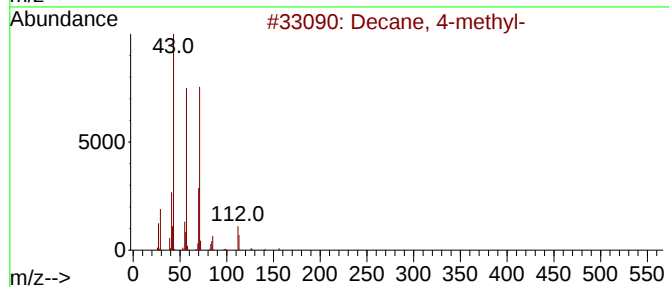
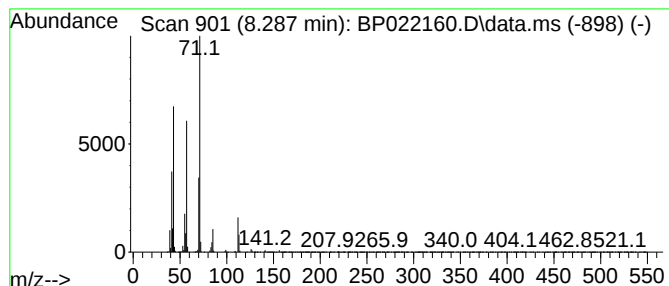
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 53

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.287	10.73 ng/ul	2312000	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 4-methyl-	156	C11H24	002847-72-5	90
2			Decane, 4-methyl-	156	C11H24	002847-72-5	76
3			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
4			Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	72
5			Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

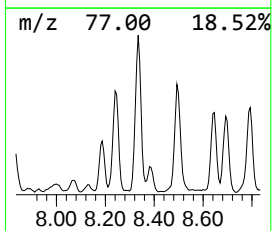
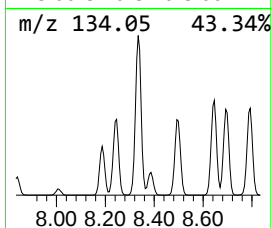
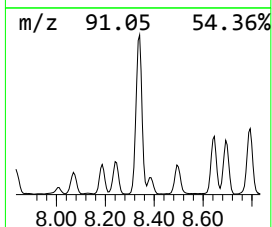
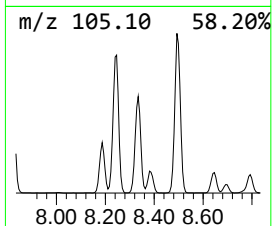
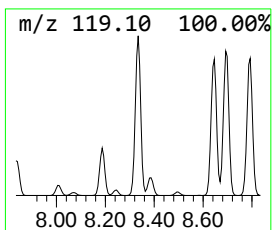
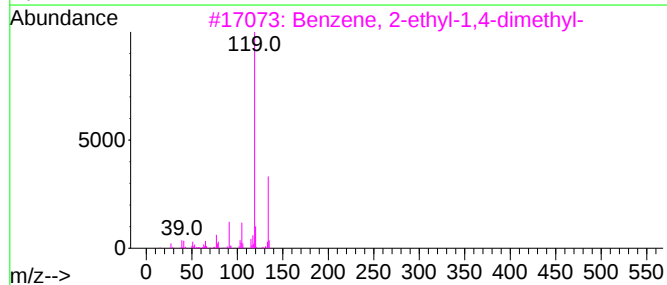
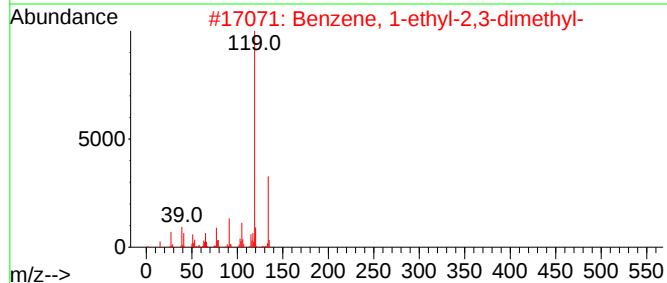
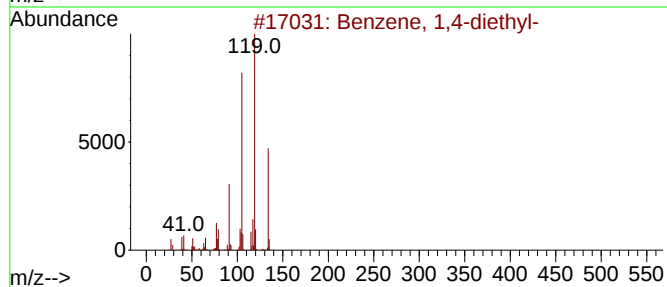
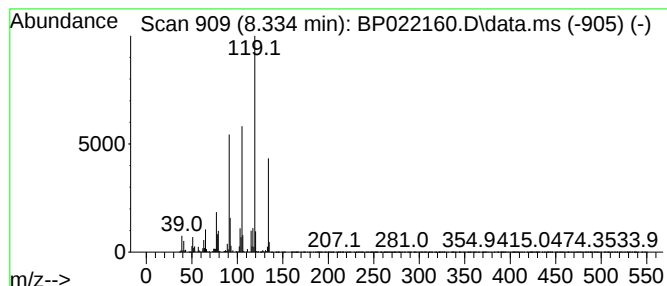
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.334	45.62 ng/ul	9824430	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	96
2			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
3			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
4			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
5			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
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Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

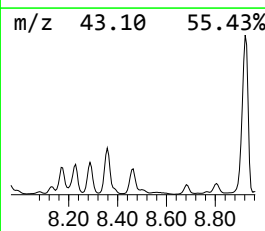
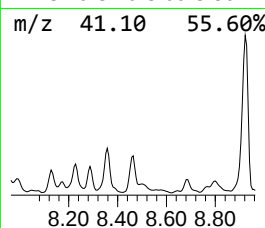
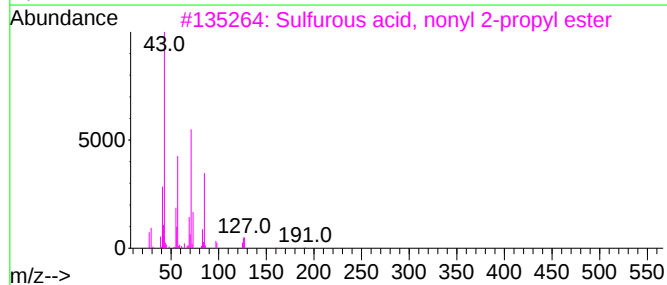
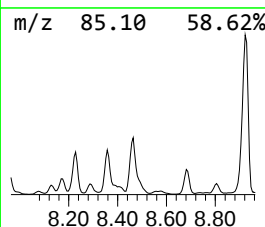
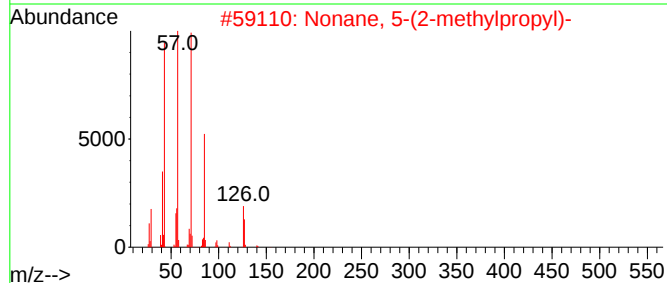
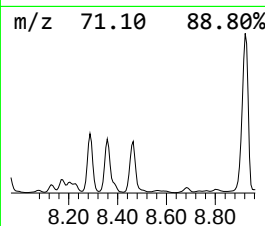
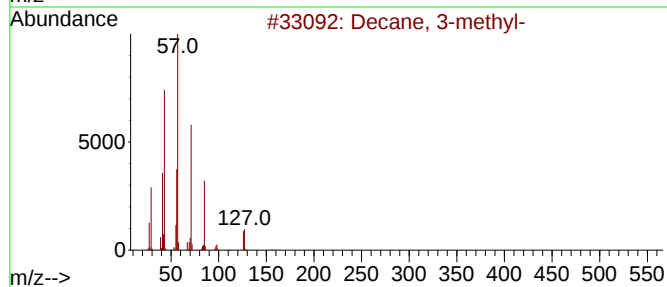
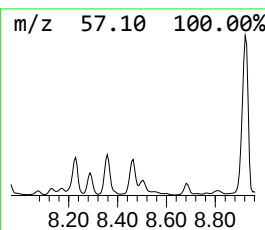
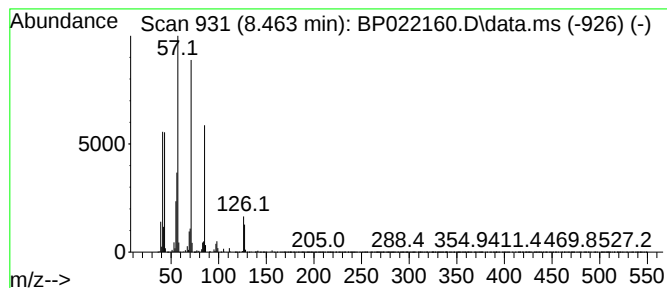
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 39

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.463	16.84 ng/ul	3627610	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane, 3-methyl-	156	C11H24	013151-34-3	91
2			Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	86
3			Sulfurous acid, nonyl 2-propyl e...	250	C12H26O3S	1000309-12-0	78
4			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	72
5			Decane, 3-methyl-	156	C11H24	013151-34-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

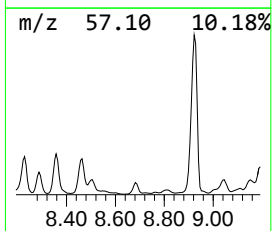
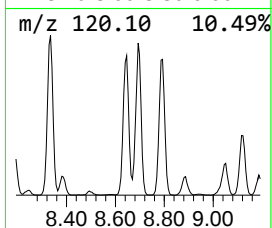
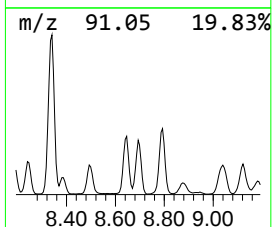
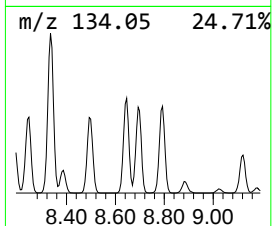
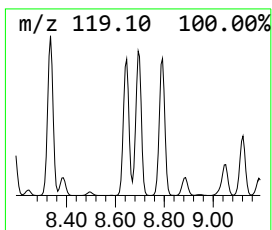
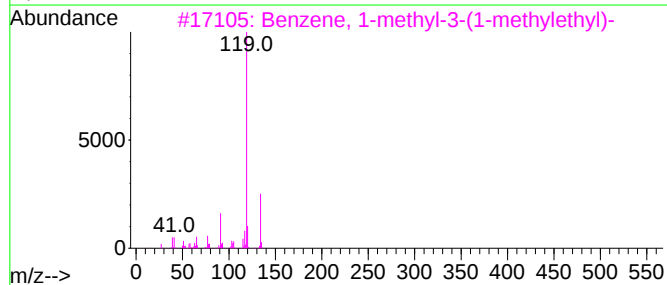
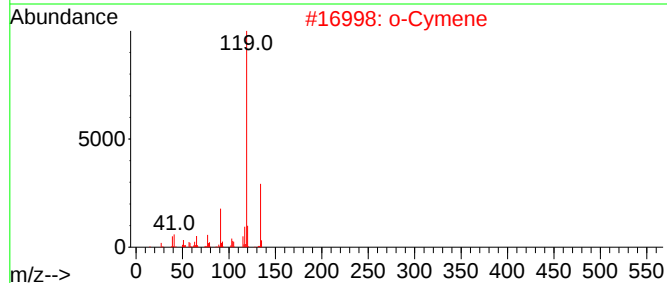
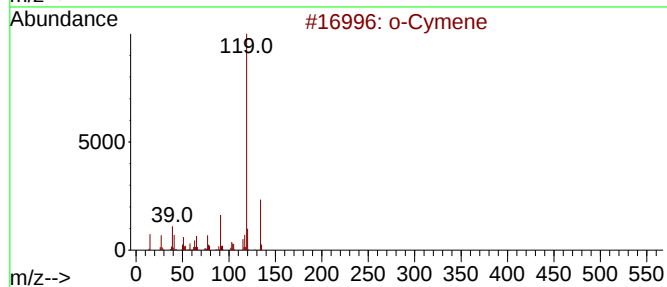
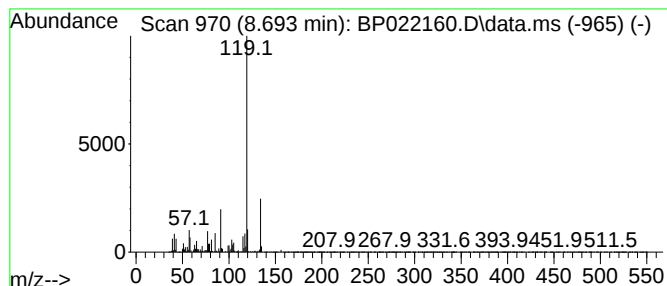
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 21 o-Cymene Concentration Rank 33

R.T.	EstConc	Area	Relative to ISTD	R.T.	
8.693	18.73 ng/ul	4033710	1,4-Dichlorobenzene-d4	7.699	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	95
2	o-Cymene	134	C10H14	000527-84-4	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	p-Cymene	134	C10H14	000099-87-6	95
5	o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

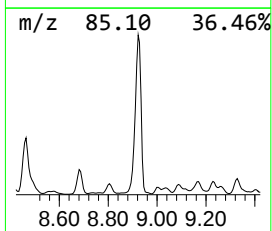
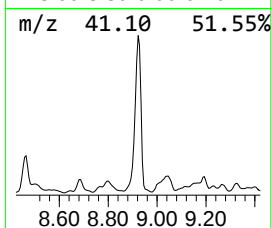
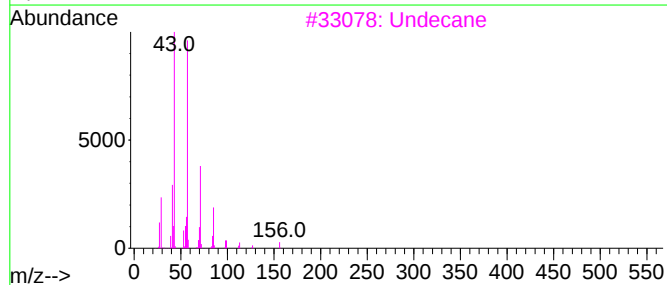
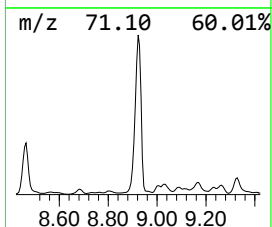
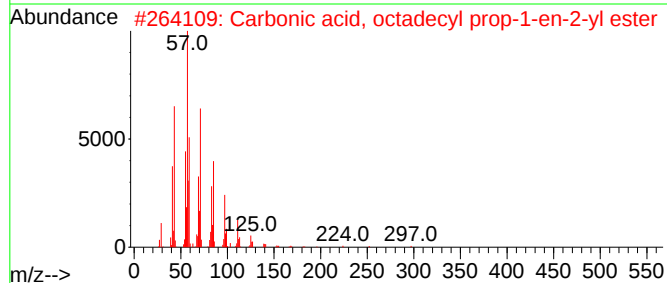
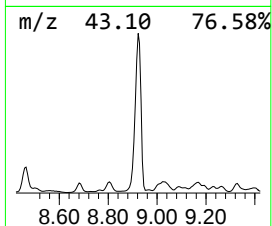
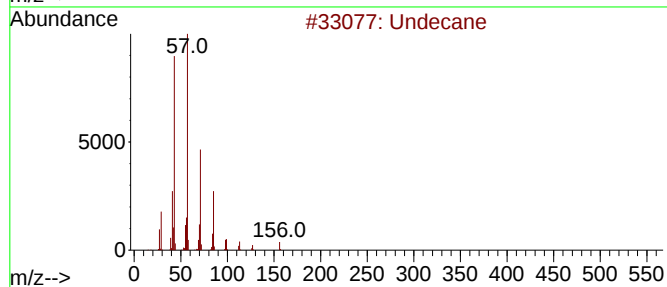
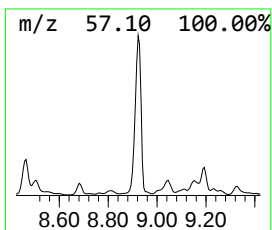
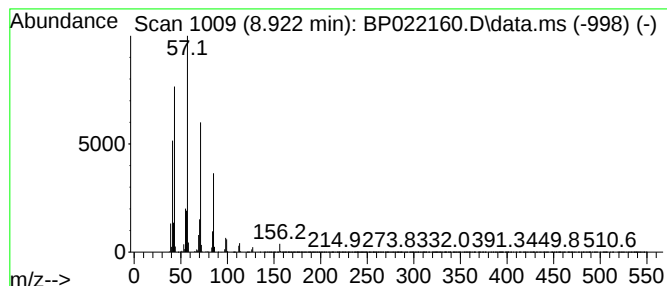
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	79.91 ng/ul	17211000	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	91
2			Carbonic acid, octadecyl prop-1-...	354	C22H42O3	1000383-11-5	90
3			Undecane	156	C11H24	001120-21-4	87
4			Tridecane	184	C13H28	000629-50-5	86
5			Hexadecane	226	C16H34	000544-76-3	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

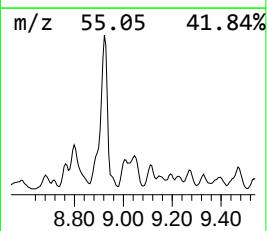
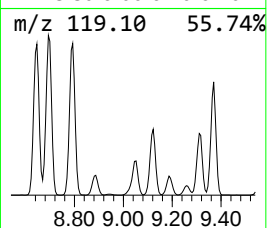
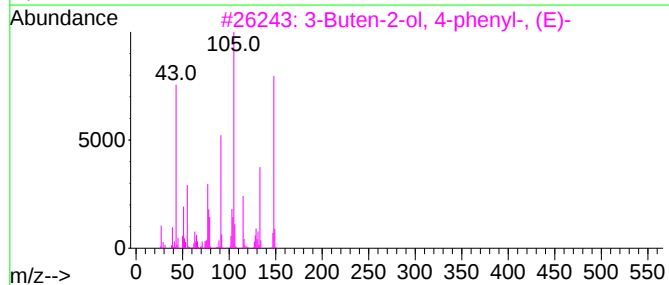
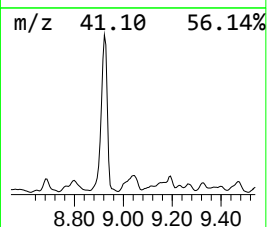
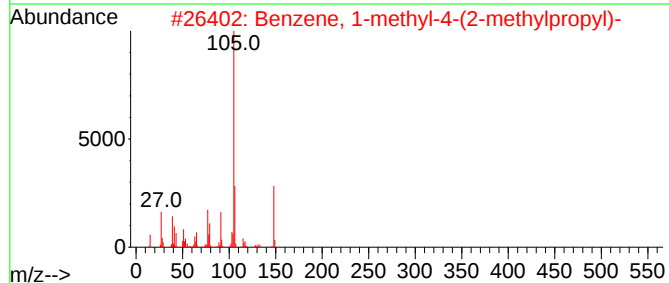
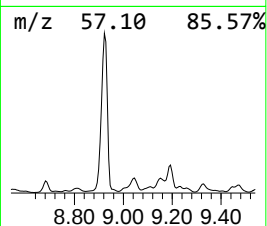
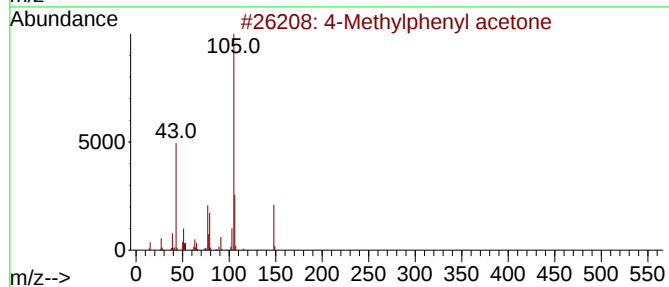
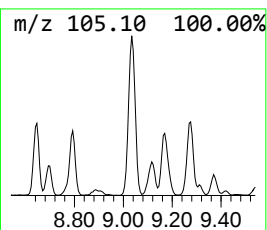
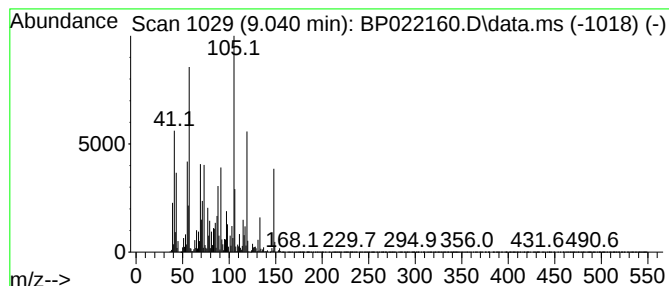
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Quant Title : SVOA CALIBRATION

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

Peak Number 23 unknown-01 Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.040	27.05 ng/ul	5825670	1,4-Dichlorobenzene-d4	7.699

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylphenyl acetone	148	C10H12O	002096-86-8	46
2			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	45
3			3-Buten-2-ol, 4-phenyl-, (E)-	148	C10H12O	1000163-70-9	38
4			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	38
5			Benzene, (1,1-dimethylpropyl)-	148	C11H16	002049-95-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

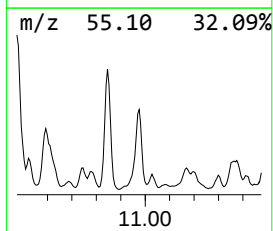
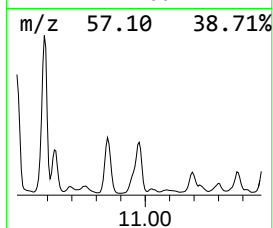
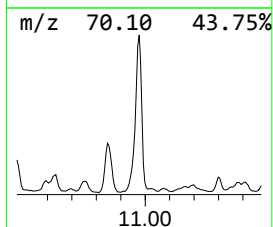
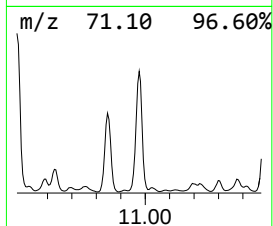
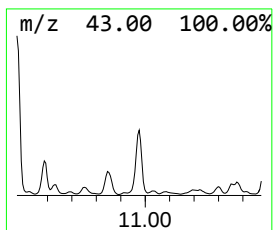
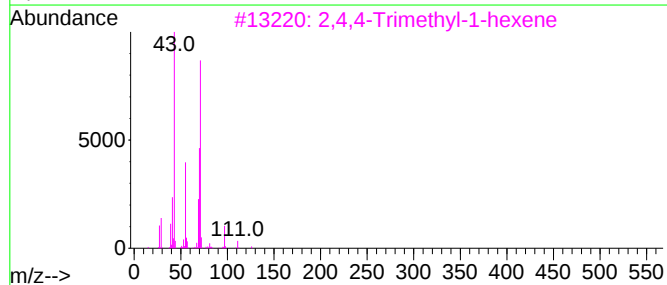
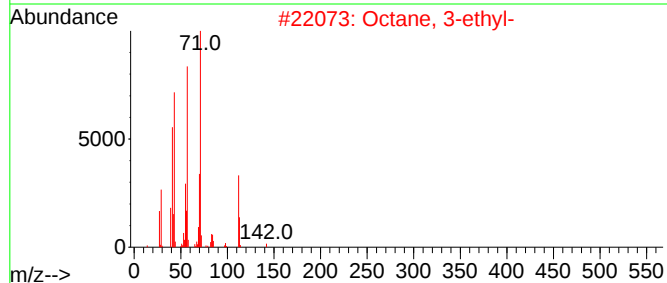
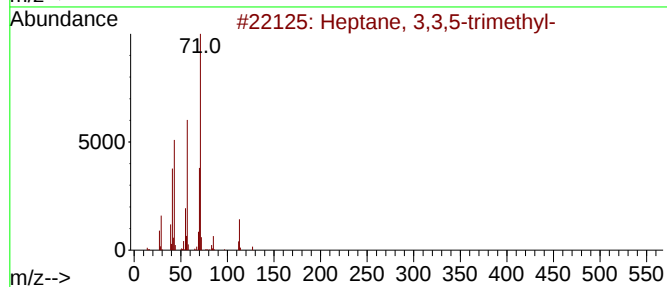
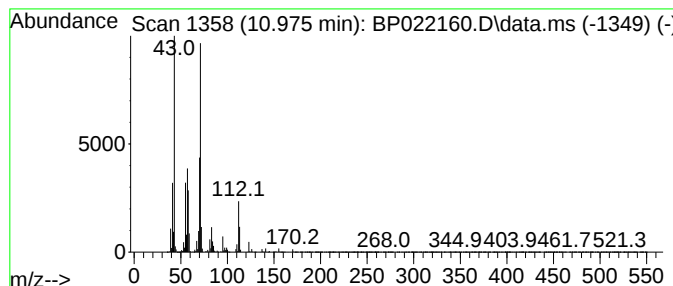
Instrument :
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ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 26 (DEL) Alkane: Straight-Chai... Concentration Rank 58

R.T.	EstConc	Area	Relative to ISTD		R.T.	
10.975	4.85 ng/ul	6954050	Naphthalene-d8		10.463	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptane, 3,3,5-trimethyl-		142	C10H22	007154-80-5	53
2	Octane, 3-ethyl-		142	C10H22	005881-17-4	53
3	2,4,4-Trimethyl-1-hexene		126	C9H18	051174-12-0	52
4	1-Butanol, 2,2-dimethyl-		102	C6H14O	001185-33-7	50
5	1-Butanol, 2,2-dimethyl-		102	C6H14O	001185-33-7	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

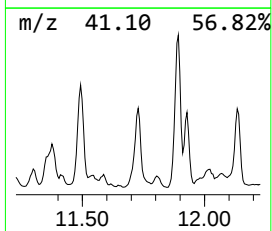
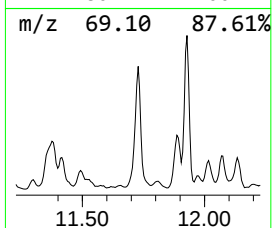
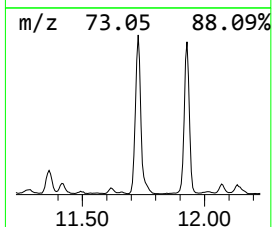
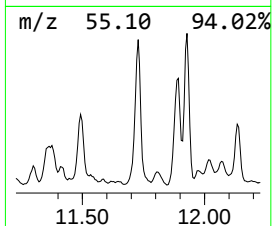
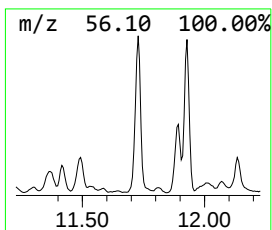
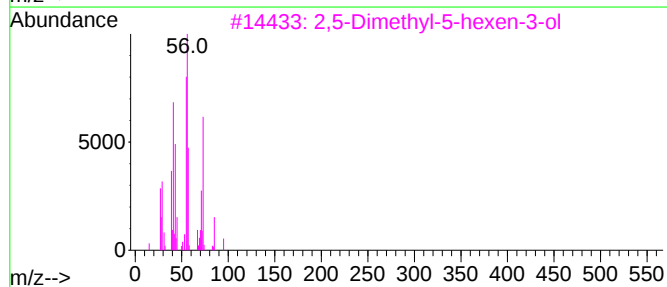
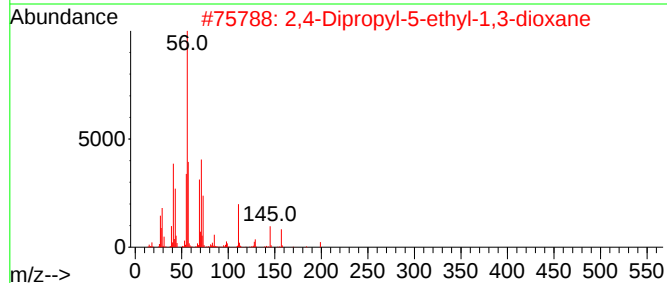
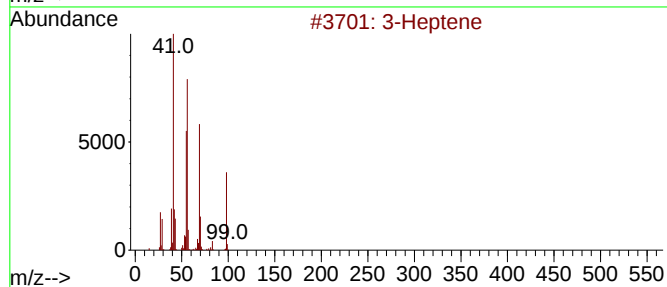
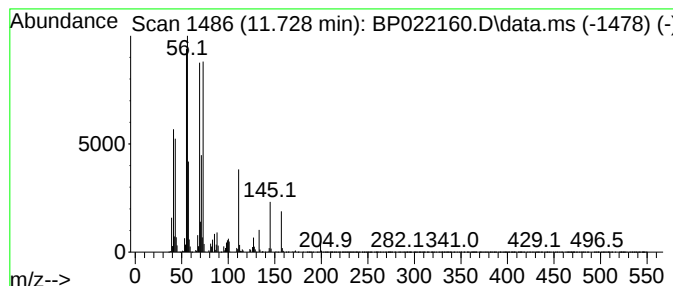
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 28 (DEL) Alkane: Straight-Chai... Concentration Rank 61

R.T.	EstConc	Area	Relative to ISTD	R.T.	
11.728	3.22 ng/ul	4612000	Naphthalene-d8	10.463	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3-Heptene	98	C7H14	000592-78-9	43
2	2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	38
3	2,5-Dimethyl-5-hexen-3-ol	128	C8H16O	067760-91-2	27
4	Cyclohexane, 1-ethyl-1,3-dimethy...	140	C10H20	062238-31-7	27
5	4-Octanol	130	C8H18O	000589-62-8	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

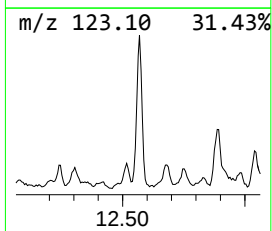
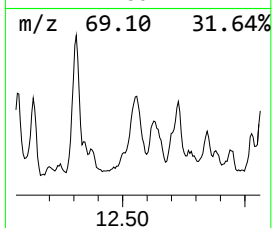
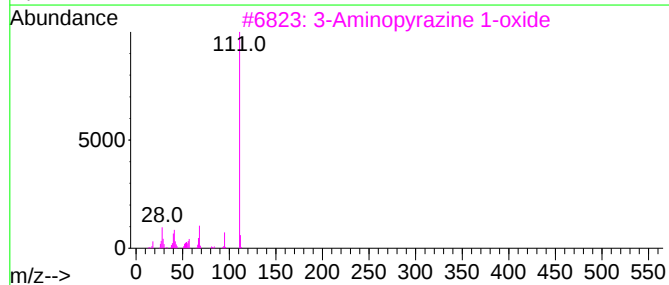
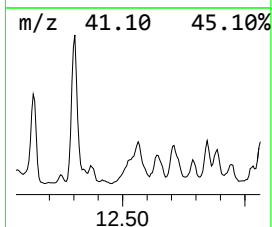
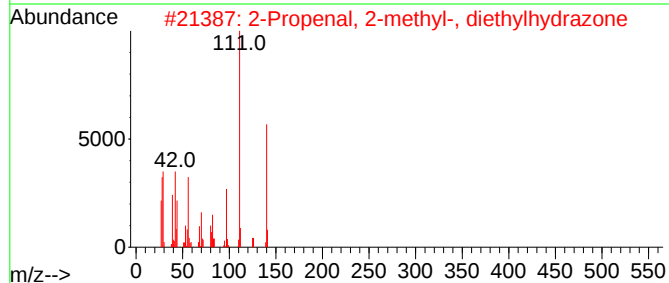
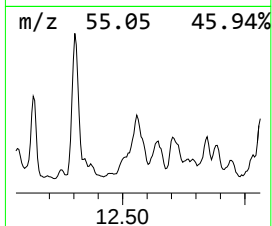
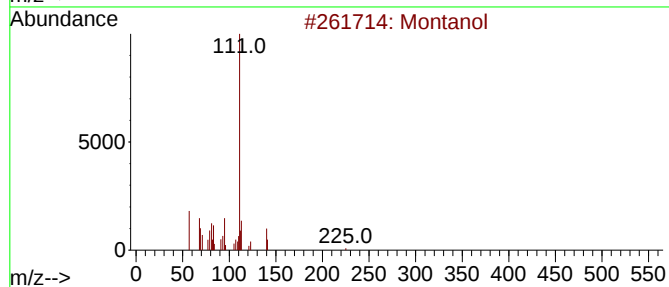
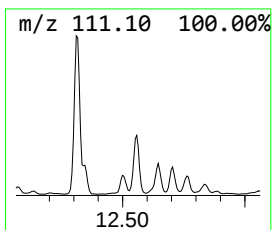
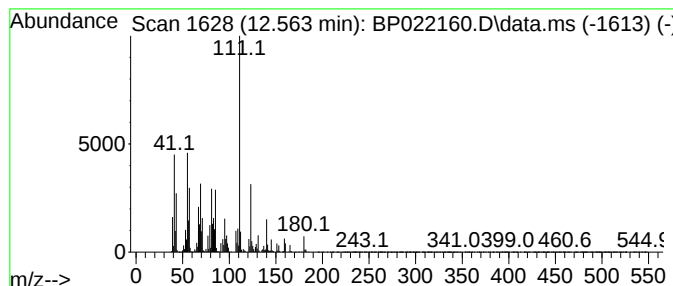
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Quant Title : SVOA CALIBRATION

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

Peak Number 31 unknown-02 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.563	36.80 ng/ul	4463230	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Montanol	352	C21H36O4	1010145-58-3	47
2			2-Propenal, 2-methyl-, diethylhy...	140	C8H16N2	075268-11-0	43
3			3-Aminopyrazine 1-oxide	111	C4H5N3O	006863-77-0	38
4			2(1H)-Pyridinone, 3-hydroxy-	111	C5H5NO2	016867-04-2	38
5			2-Aminopyrimidine-1-oxide	111	C4H5N3O	035034-15-2	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

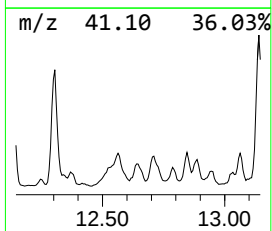
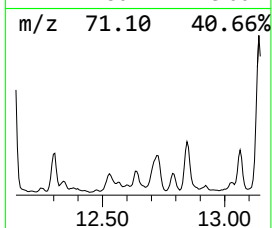
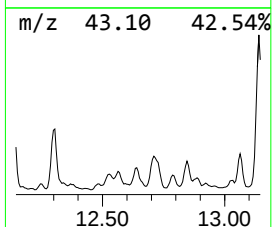
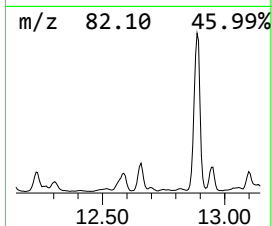
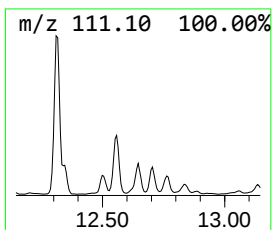
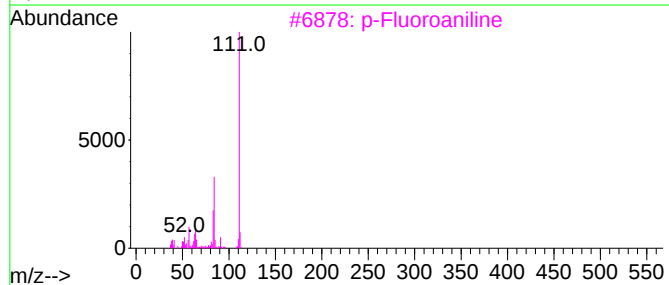
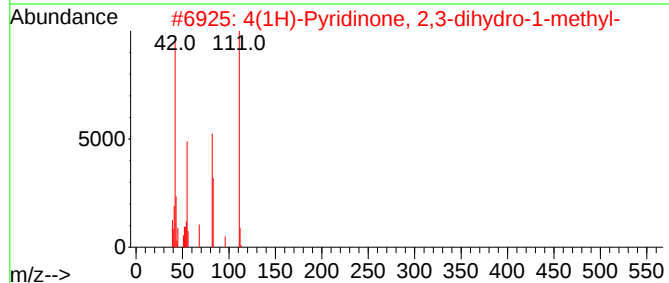
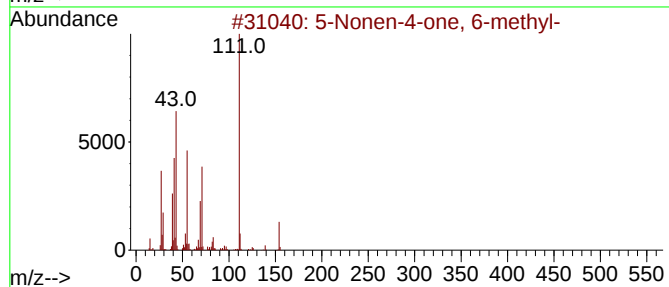
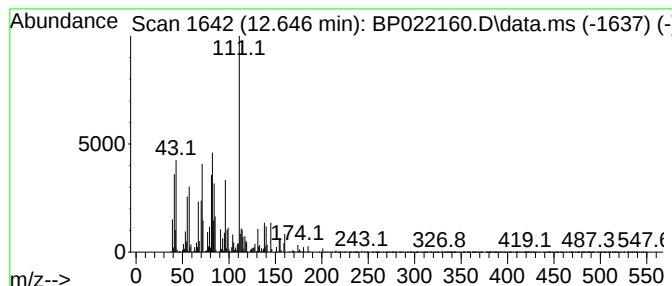
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Quant Title : SVOA CALIBRATION

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

Peak Number 32 unknown-03 Concentration Rank 37

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.646	17.04 ng/ul	2067220	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Nonen-4-one, 6-methyl-	154	C10H18O	007036-98-8	35
2			4(1H)-Pyridinone, 2,3-dihydro-1-...	111	C6H9NO	035488-00-7	32
3			p-Fluoroaniline	111	C6H6FN	000371-40-4	30
4			Benzenamine, 3-fluoro-	111	C6H6FN	000372-19-0	27
5			Pentanal, 2-methylene-, 2-(1-met...	154	C9H18N2	033063-80-8	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

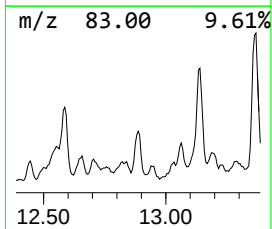
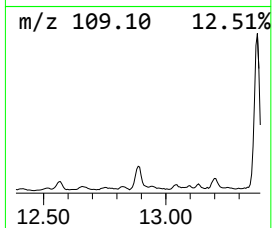
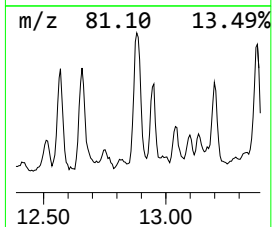
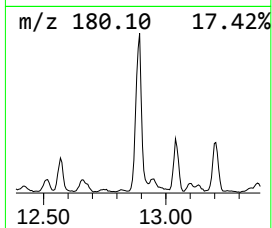
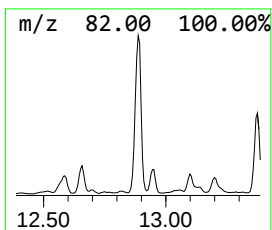
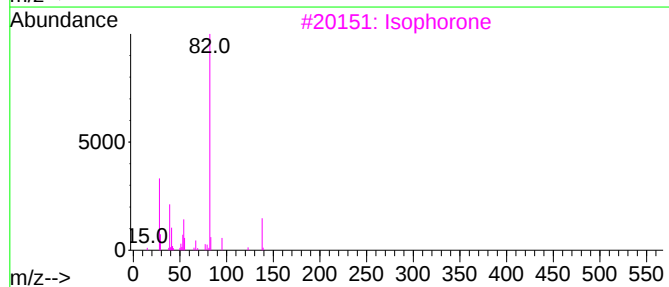
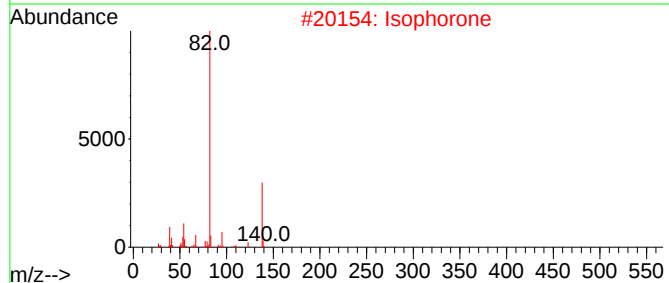
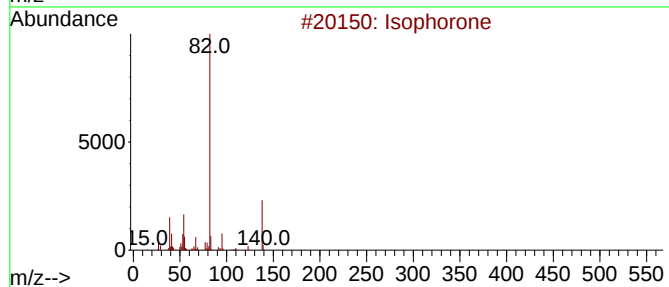
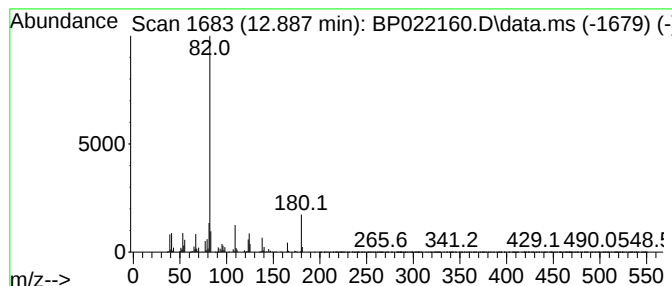
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BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 33 1H-Pyrazole, 4,5-dihydro-5,... Concentration Rank 38

R.T.	EstConc	Area	Relative to ISTD		R.T.	
12.887	17.01 ng/ul	2063460	Acenaphthene-d10		14.328	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Isophorone		138	C9H14O	000078-59-1	53
2	Isophorone		138	C9H14O	000078-59-1	53
3	Isophorone		138	C9H14O	000078-59-1	53
4	1H-Pyrazole, 4,5-dihydro-5,5-dim...		138	C8H14N2	106251-09-6	53
5	cis-Hexenyl octyne carbonate		236	C15H24O2	068480-29-5	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

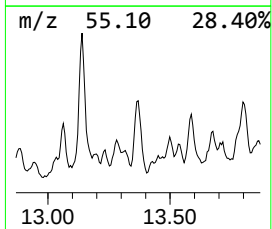
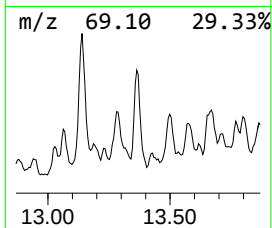
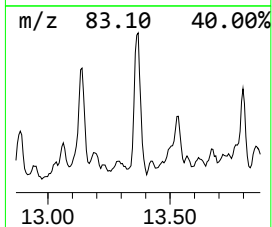
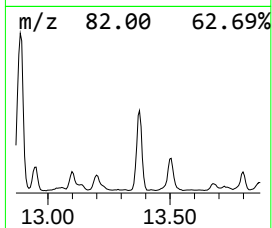
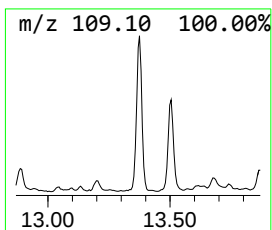
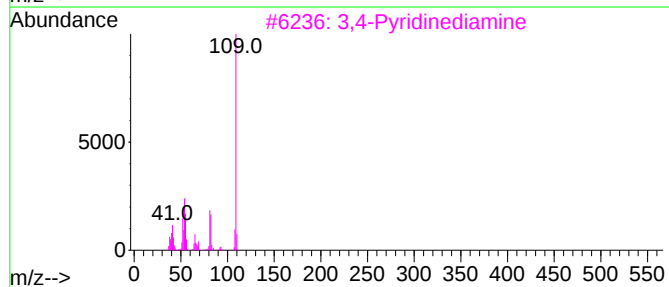
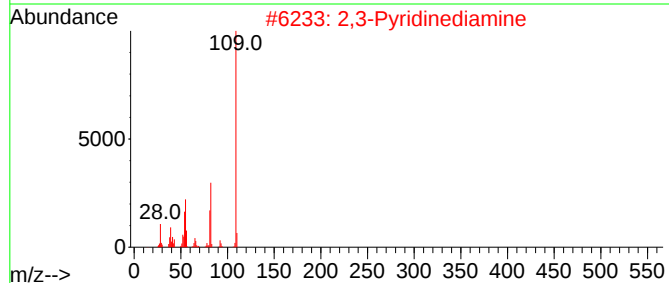
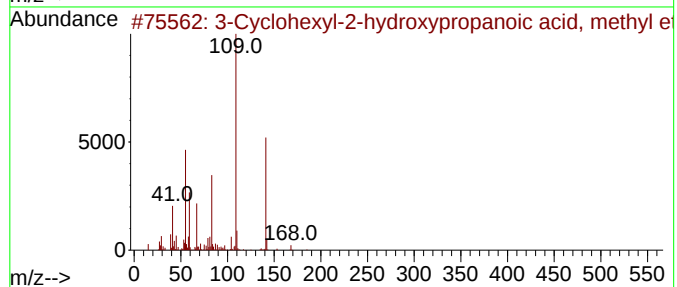
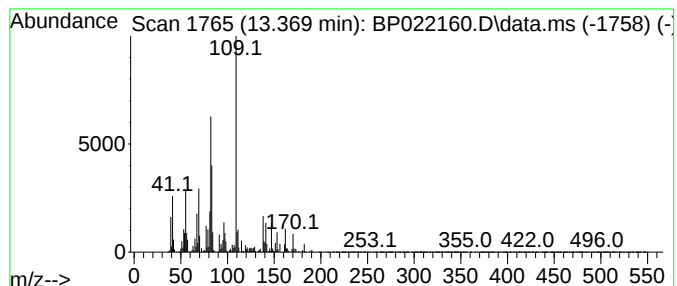
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Quant Title : SVOA CALIBRATION

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

Peak Number 34 unknown-04 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.369	47.81 ng/ul	5798210	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3-Cyclohexyl-2-hydroxypropanoic ...	200	C11H20O3	078811-34-4	47
2			2,3-Pyridinediamine	109	C5H7N3	000452-58-4	43
3			3,4-Pyridinediamine	109	C5H7N3	000054-96-6	30
4			3-Heptyne, 5-ethyl-5-methyl-	138	C10H18	061228-10-2	30
5			Cyclopentane, (2-methylbutylidene)-	138	C10H18	053366-54-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

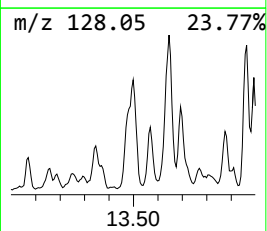
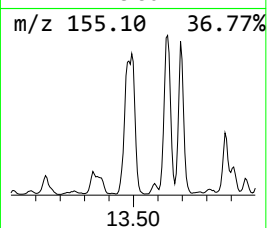
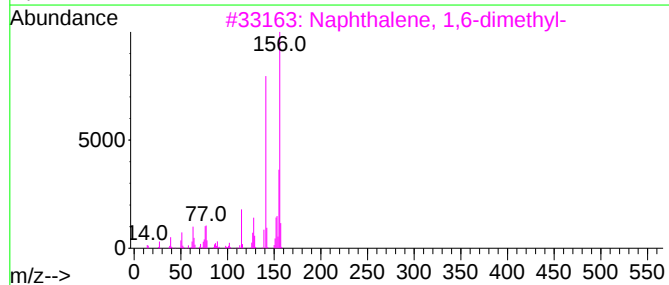
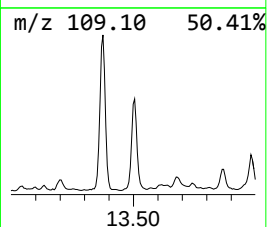
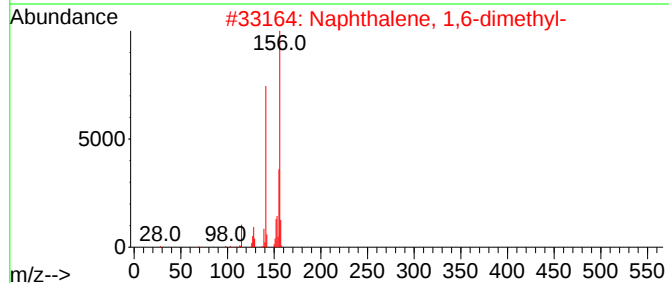
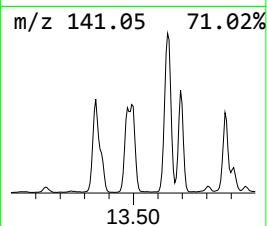
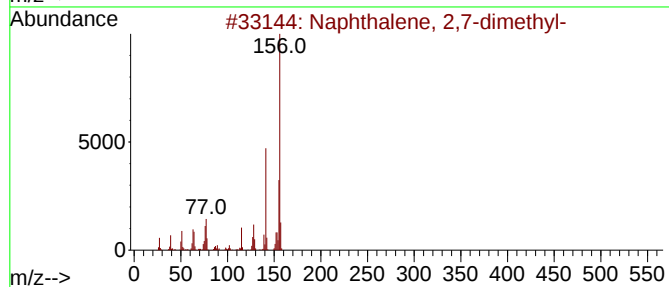
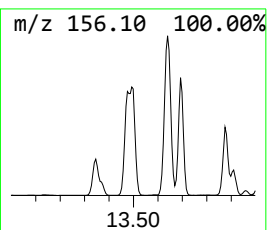
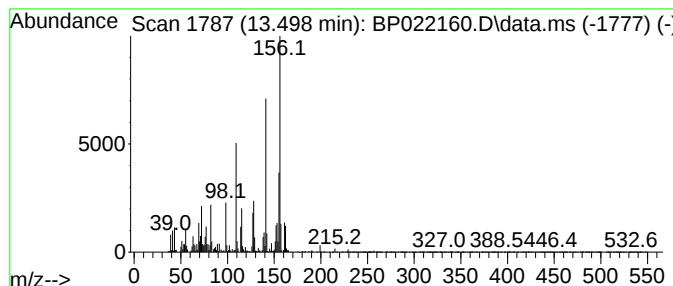
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Quant Title : SVOA CALIBRATION

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

Peak Number 35 Naphthalene, 2,7-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.499	54.20 ng/ul	6573840	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
2			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	90
4			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90
5			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

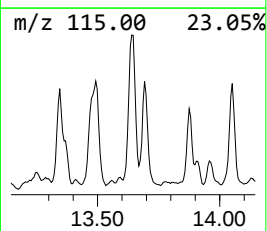
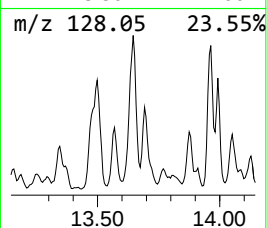
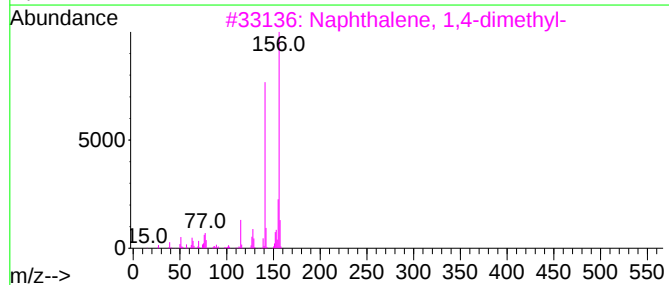
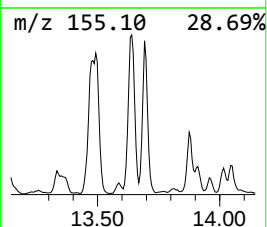
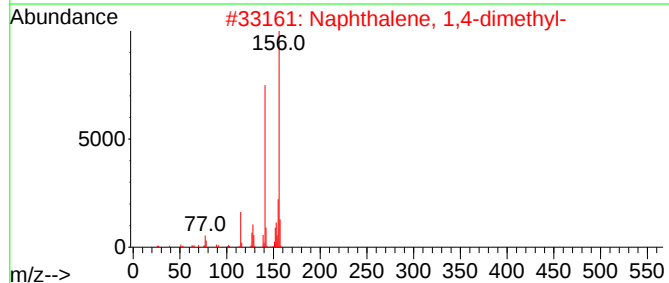
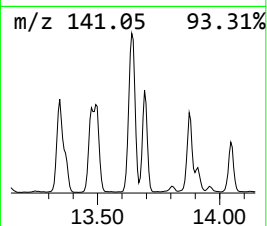
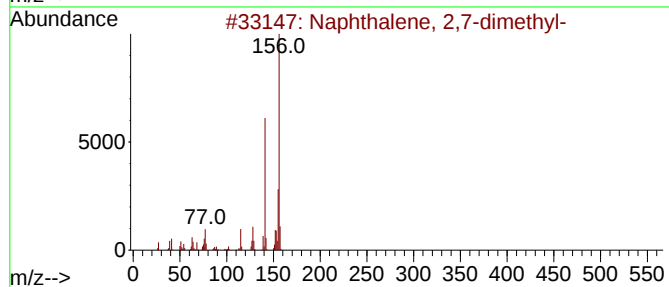
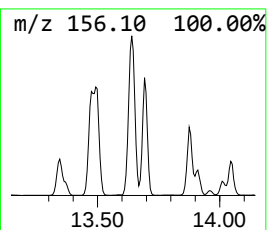
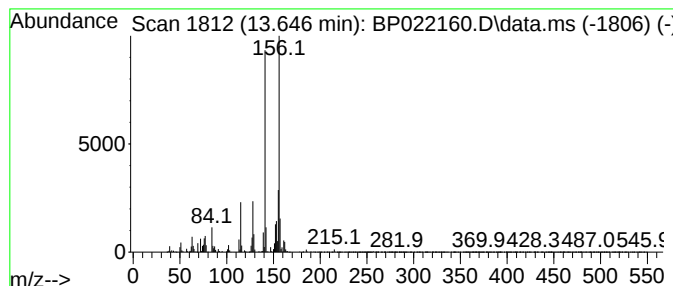
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Quant Title : SVOA CALIBRATION

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

Peak Number 36 Naphthalene, 1,4-dimethyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.646	35.87 ng/ul	4349970	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
3			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	95
4			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	95
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

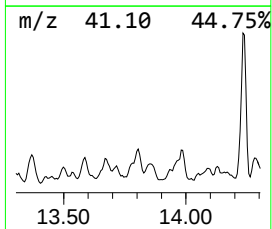
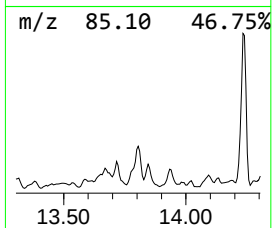
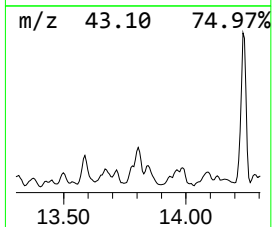
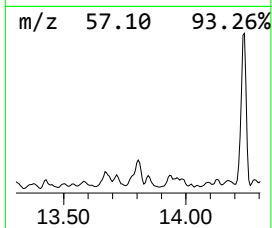
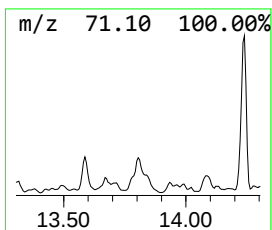
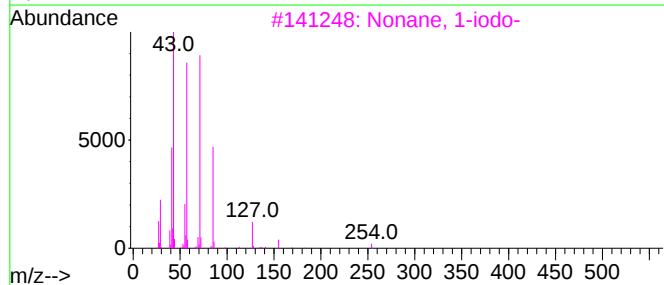
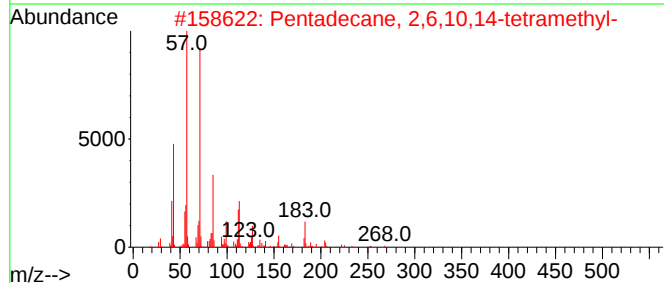
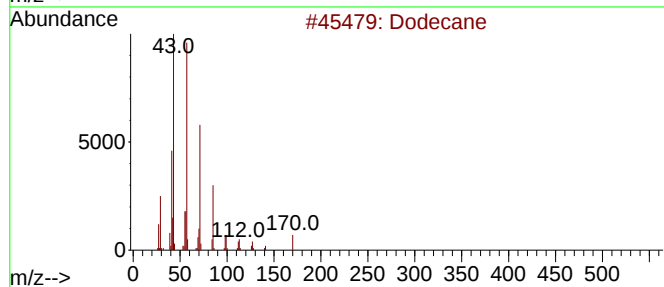
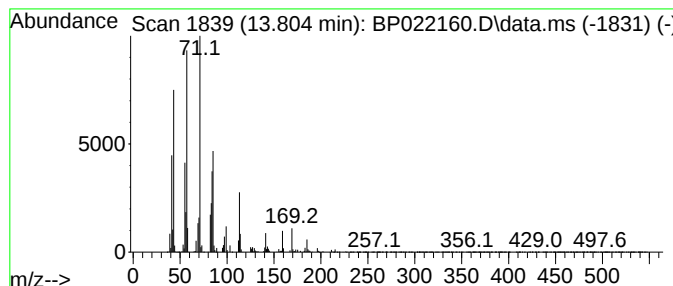
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Quant Title : SVOA CALIBRATION

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

Peak Number 37 (DEL) Alkane: Straight-Chai... Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.804	25.50 ng/ul	3092450	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dodecane	170	C12H26	000112-40-3	64
2			Pentadecane, 2,6,10,14-tetramethyl-	268	C19H40	001921-70-6	53
3			Nonane, 1-iodo-	254	C9H19I	004282-42-2	47
4			Undecane, 5-methyl-	170	C12H26	001632-70-8	46
5			Carbonic acid, dodecyl vinyl ester	256	C15H28O3	1000382-54-8	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

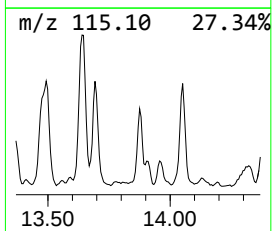
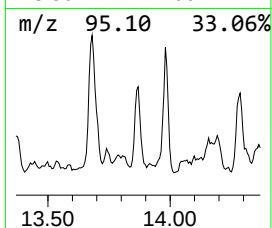
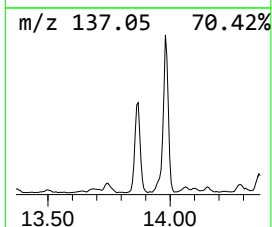
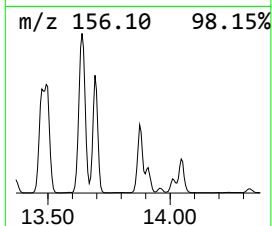
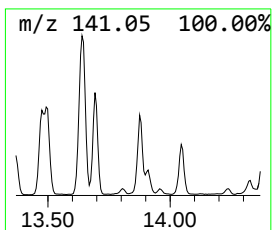
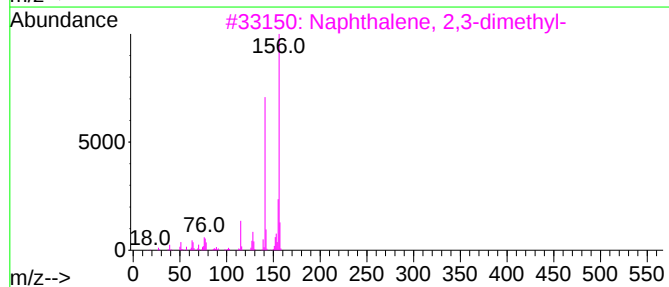
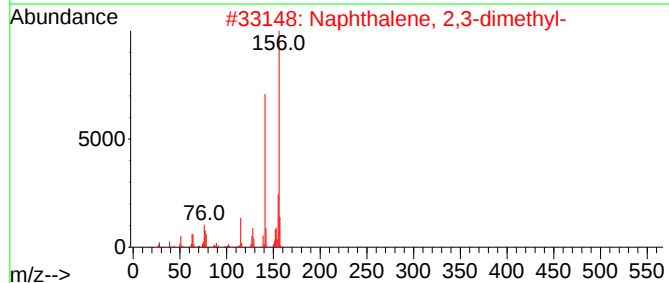
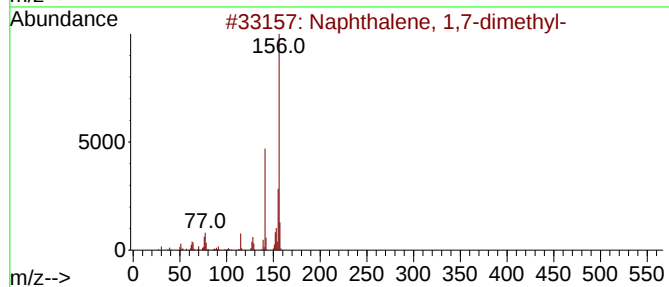
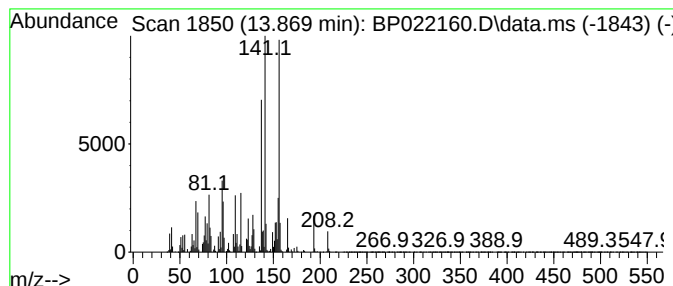
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 38 Naphthalene, 1,7-dimethyl- Concentration Rank 26

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.869	21.45 ng/ul	2601780	Acenaphthene-d10	14.328	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
2	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	91
3	Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	91
4	Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	91
5	Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
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Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

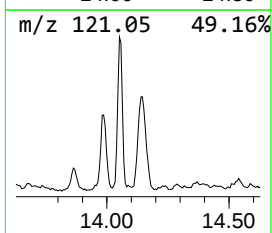
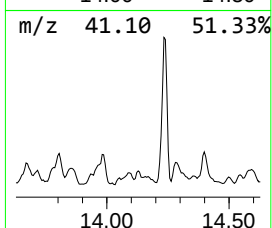
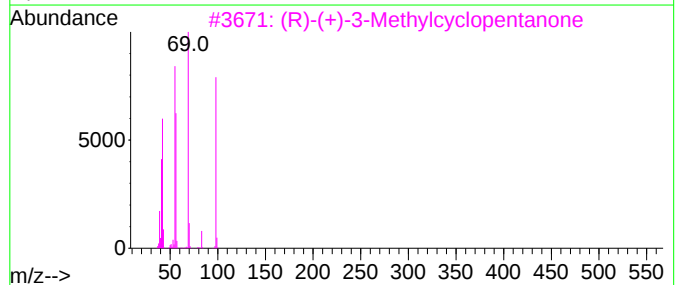
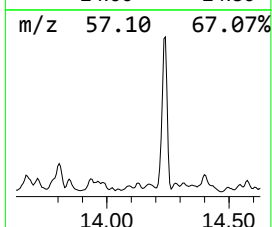
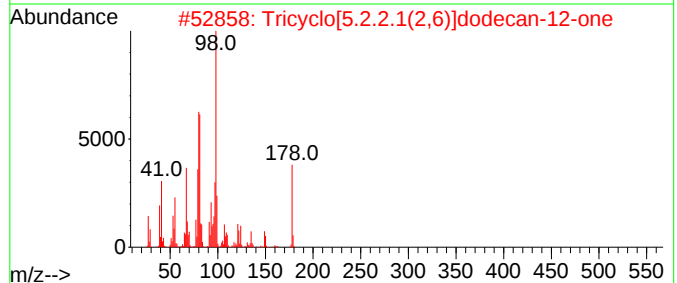
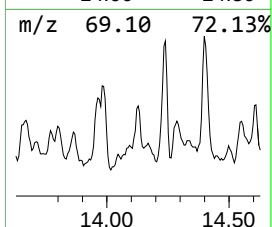
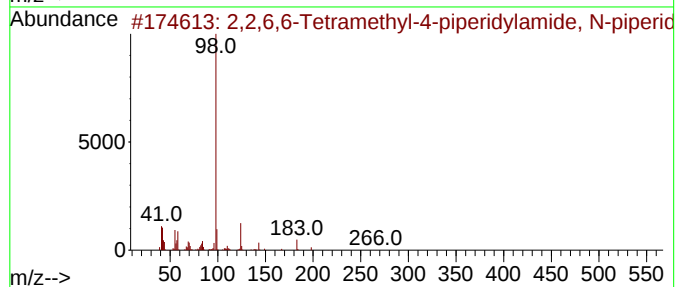
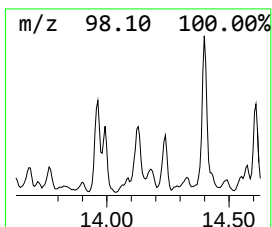
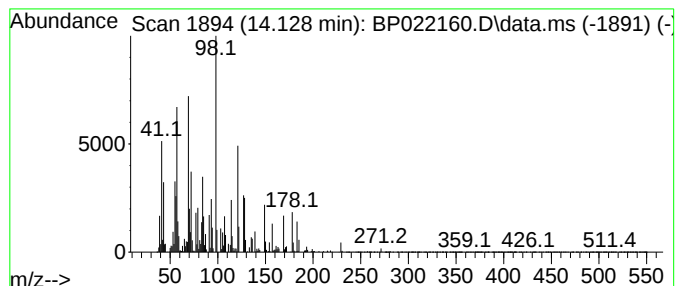
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Quant Title : SVOA CALIBRATION

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

Peak Number 39 unknown-05 Concentration Rank 28

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.128	20.23 ng/ul	2453650	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2,6,6-Tetramethyl-4-piperidyla...	281	C16H31N3O	109172-31-8	30		
2	Tricyclo[5.2.2.1(2,6)]dodecan-12...	178	C12H18O	1000185-61-9	27		
3	(R)-(+)-3-Methylcyclopentanone	98	C6H10O	006672-30-6	25		
4	Methyl 5-piperidino-4-ketocaproate	213	C11H19NO3	018904-36-4	25		
5	Cyclohexanone, 2-butyl-	154	C10H18O	001126-18-7	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

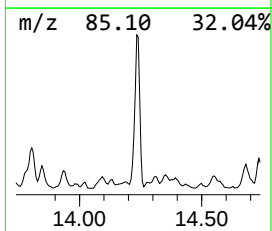
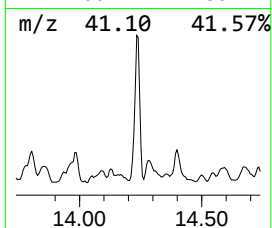
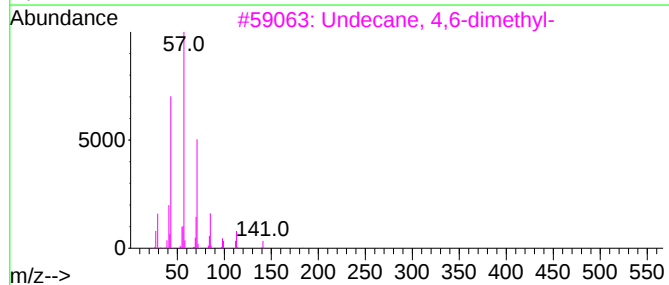
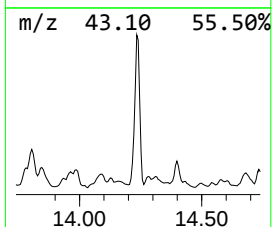
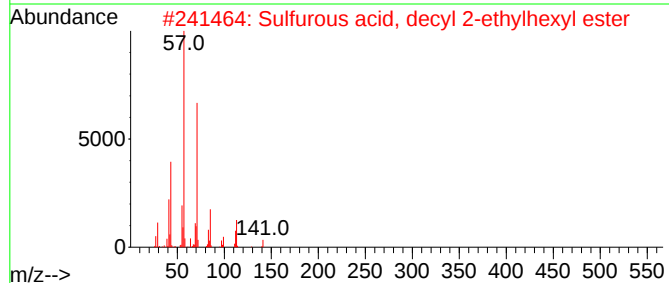
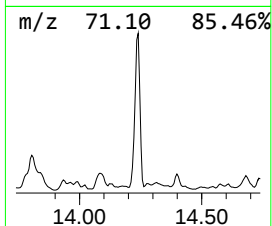
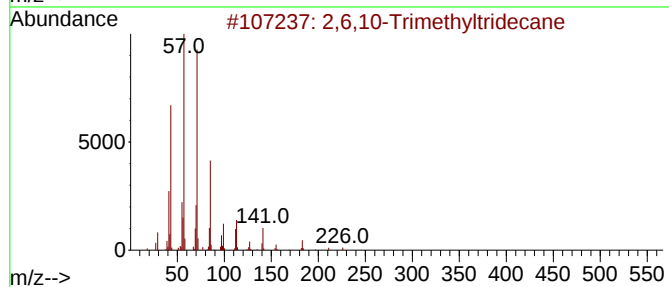
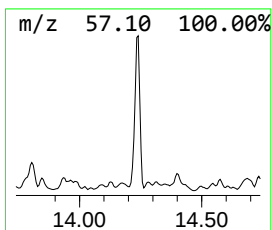
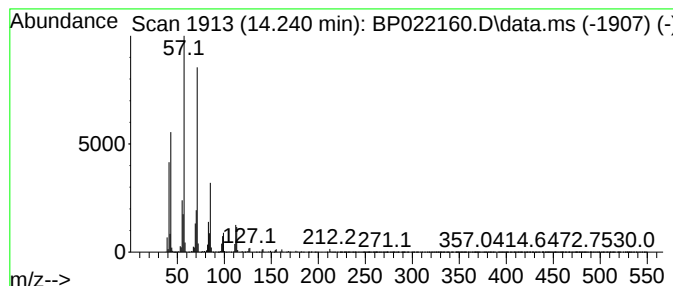
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Quant Title : SVOA CALIBRATION

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

Peak Number 40 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.240	70.48 ng/ul	8548470	Acenaphthene-d10	14.328

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,6,10-Trimethyltridecane	226	C16H34	003891-99-4	86	
2	Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	86	
3	Undecane, 4,6-dimethyl-	184	C13H28	017312-82-2	86	
4	Undecane, 2,6-dimethyl-	184	C13H28	017301-23-4	80	
5	Decane, 2-methyl-	156	C11H24	006975-98-0	72	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

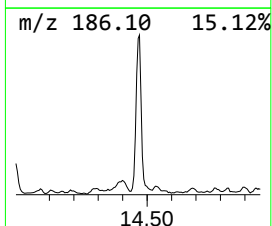
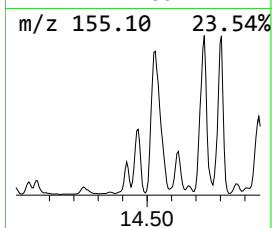
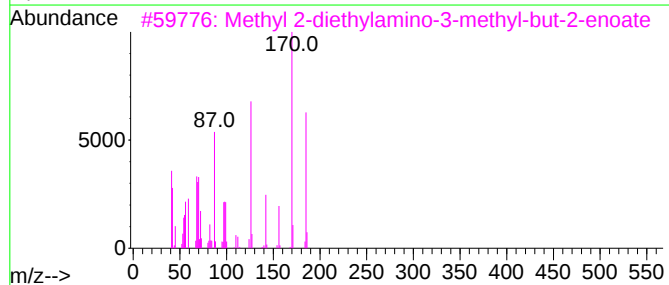
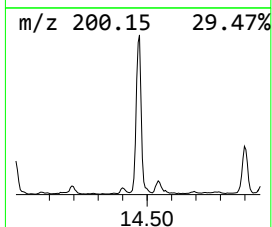
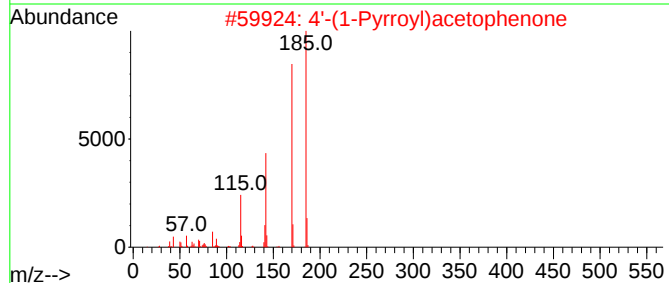
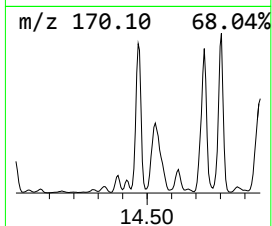
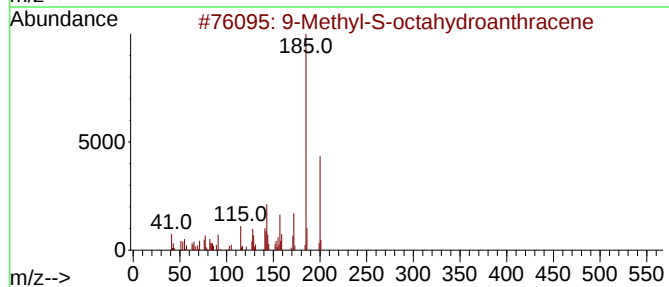
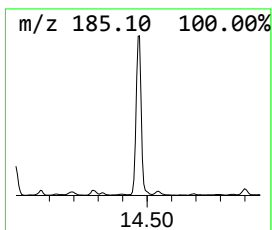
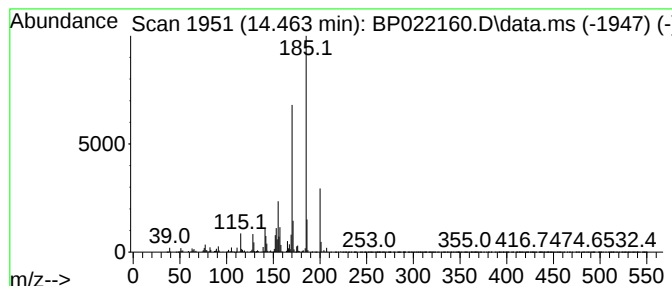
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Quant Title : SVOA CALIBRATION

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

Peak Number 41 9-Methyl-S-octahydroanthracene Concentration Rank 25

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.463	21.89 ng/ul	2654400	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	64
2			4'-(1-Pyrrolyl)acetophenone	185	C12H11NO	022106-37-2	50
3			Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	47
4			1H-Pyrazole, 5-(t-butyl)-3-phenyl-	200	C13H16N2	1000261-34-8	46
5			6-Fluoro-2,4-pyrimidinediamine, TMS	200	C7H13FN4Si	1000472-66-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

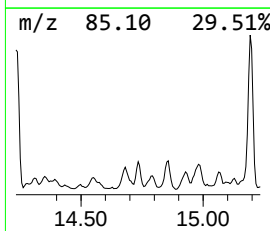
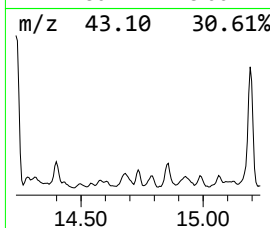
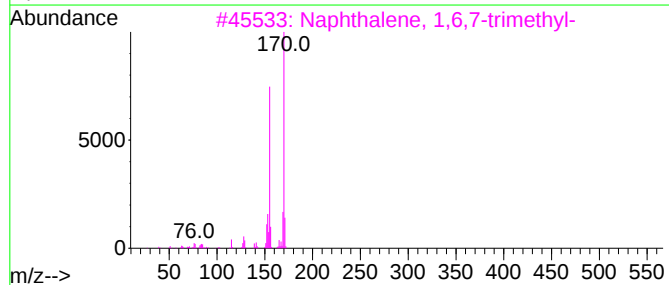
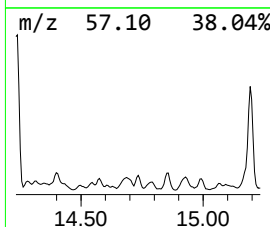
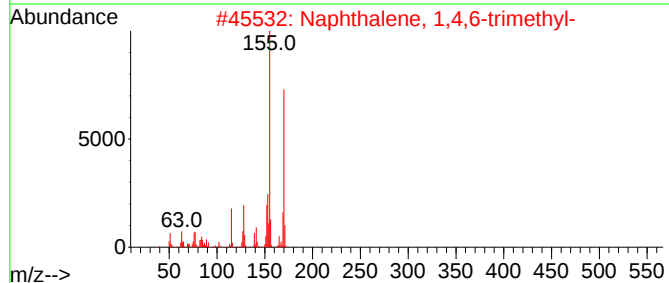
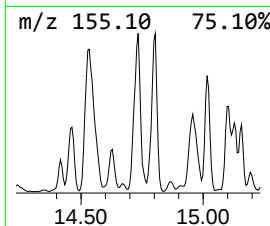
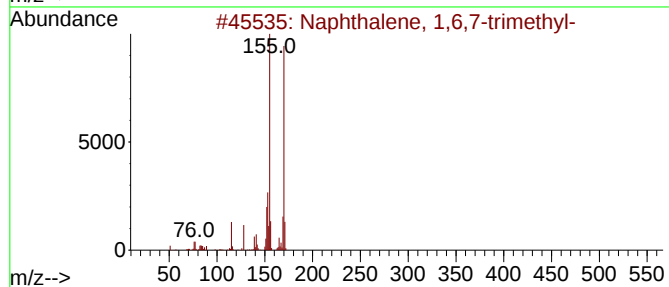
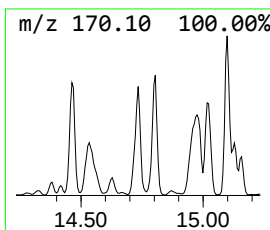
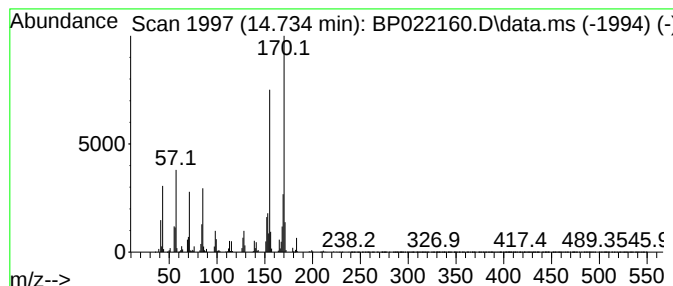
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Quant Title : SVOA CALIBRATION

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

Peak Number 42 Naphthalene, 1,6,7-trimethyl- Concentration Rank 23

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.734	22.94 ng/ul	2782520	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	93
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	93
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	91
4			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	91
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

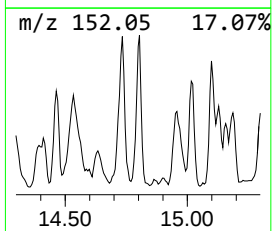
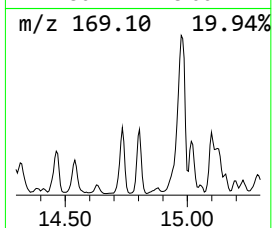
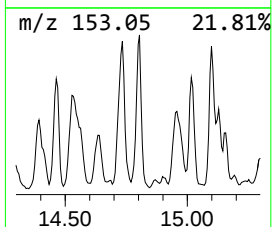
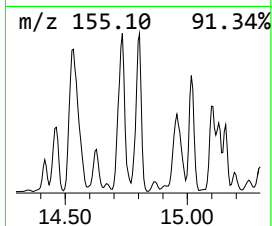
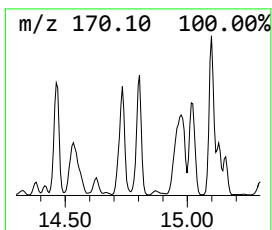
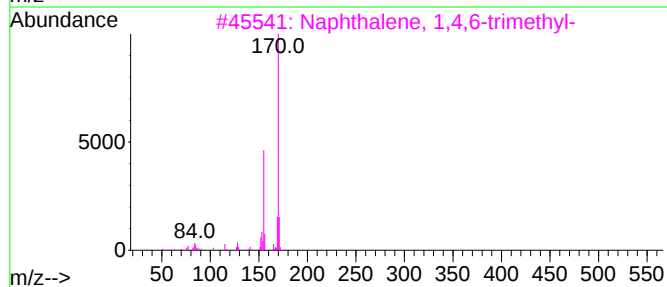
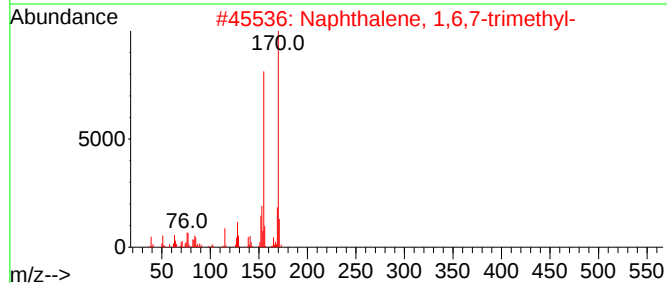
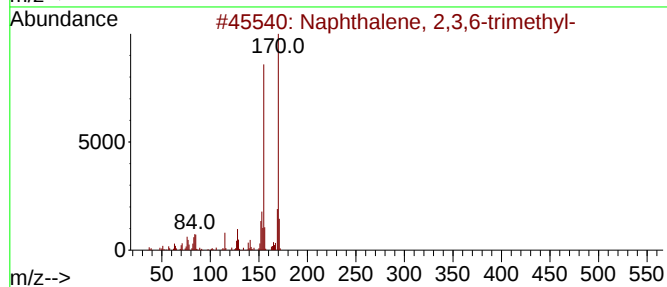
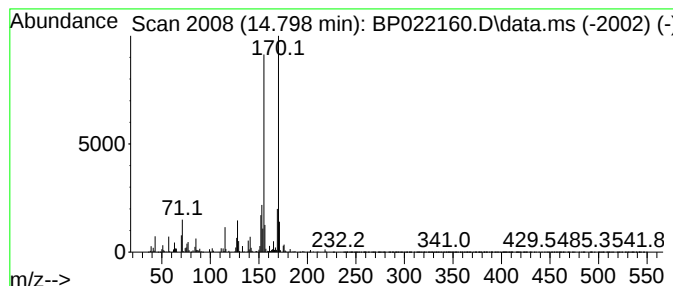
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Quant Title : SVOA CALIBRATION

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

Peak Number 43 Naphthalene, 2,3,6-trimethyl- Concentration Rank 27

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.798	21.21 ng/ul	2572510	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	97
2			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	95
3			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	94
4			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	94
5			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
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Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

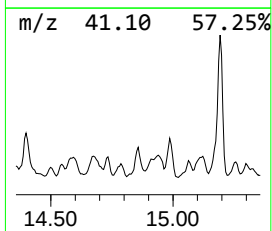
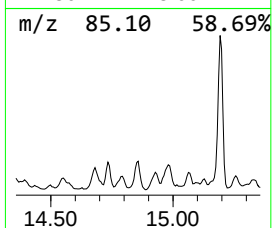
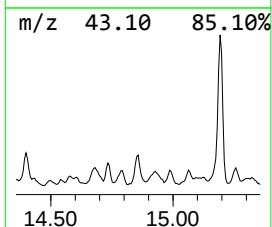
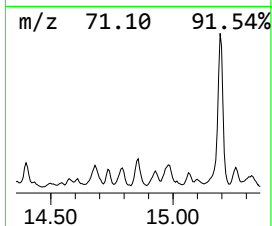
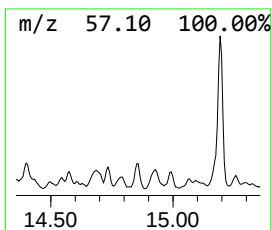
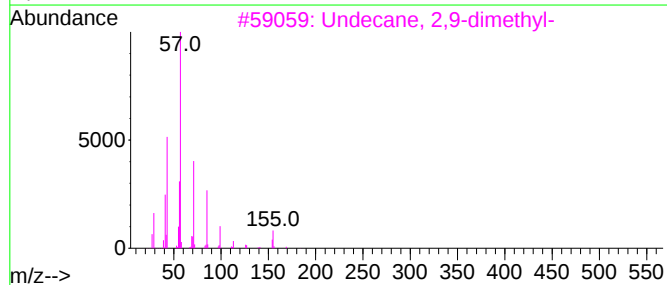
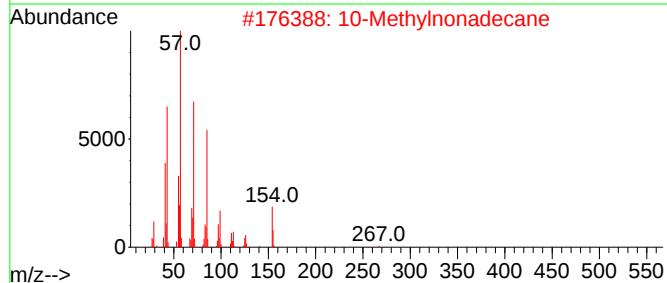
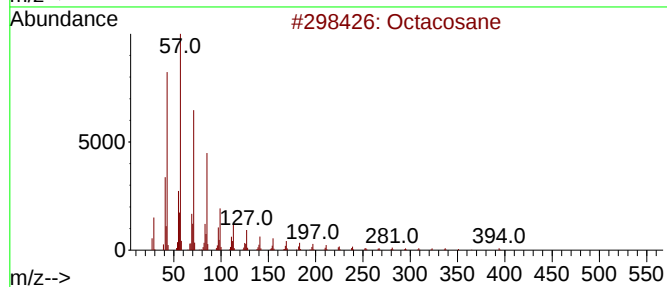
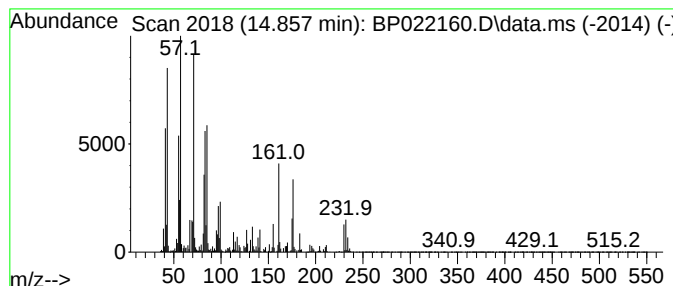
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Quant Title : SVOA CALIBRATION

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

Peak Number 44 (DEL) Alkane: Straight-Chai... Concentration Rank 24

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.857	22.64 ng/ul	2746070	Acenaphthene-d10	14.328	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octacosane	394	C28H58	000630-02-4	46
2	10-Methylnonadecane	282	C20H42	056862-62-5	46
3	Undecane, 2,9-dimethyl-	184	C13H28	017301-26-7	43
4	Tridecane, 2-methyl-	198	C14H30	001560-96-9	43
5	Pentacosane	352	C25H52	000629-99-2	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
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Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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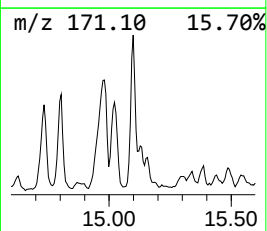
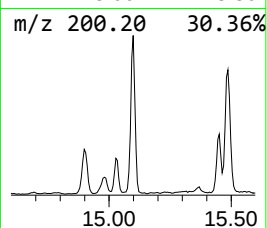
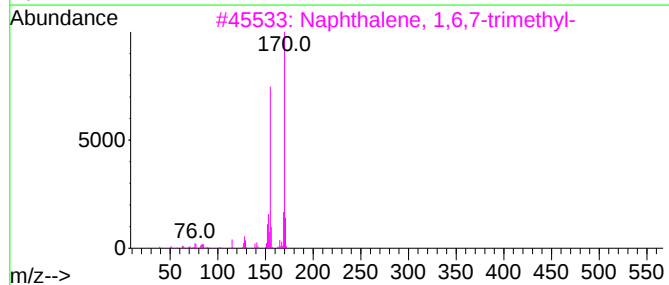
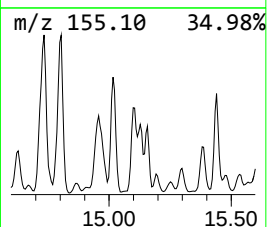
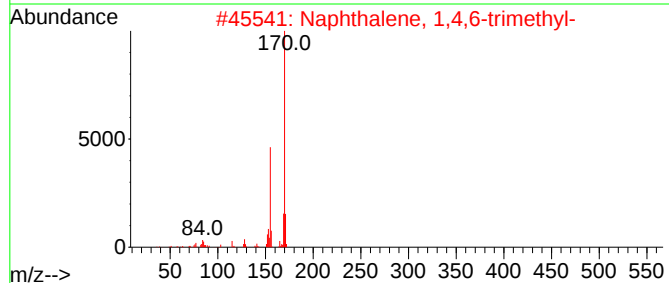
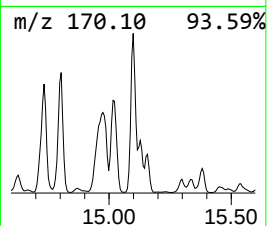
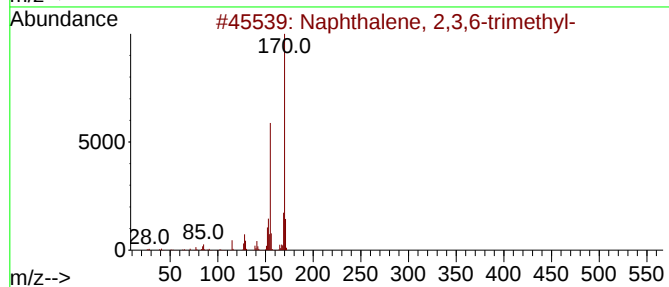
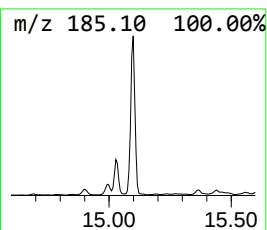
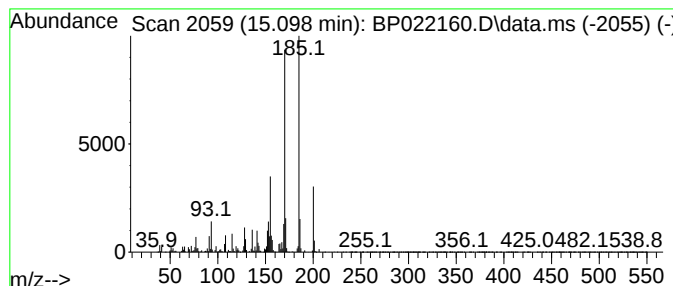
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Quant Title : SVOA CALIBRATION

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

Peak Number 45 Naphthalene, 1,4,6-trimethyl- Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.098	30.32 ng/ul	3677840	Acenaphthene-d10	14.328

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	90
2			Naphthalene, 1,4,6-trimethyl-	170	C13H14	002131-42-2	89
3			Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	64
4			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55
5			4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

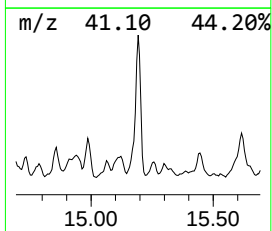
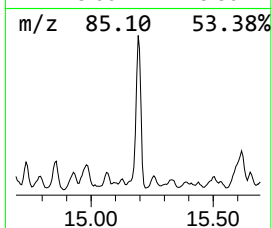
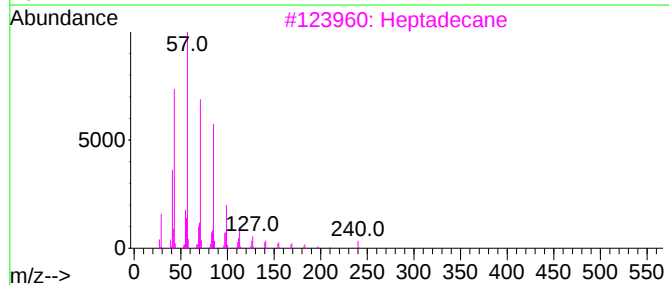
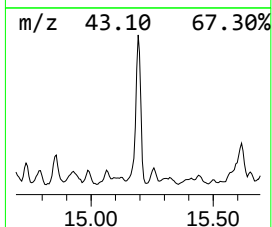
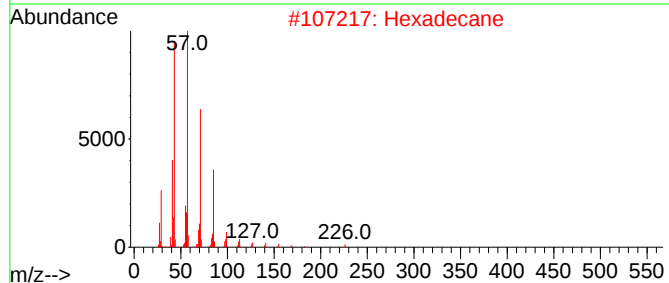
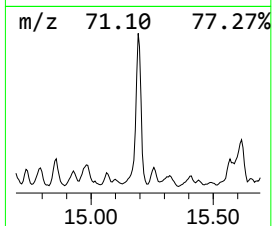
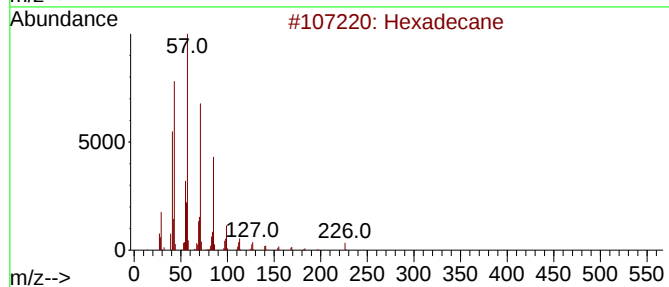
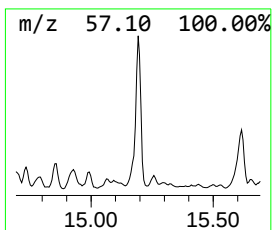
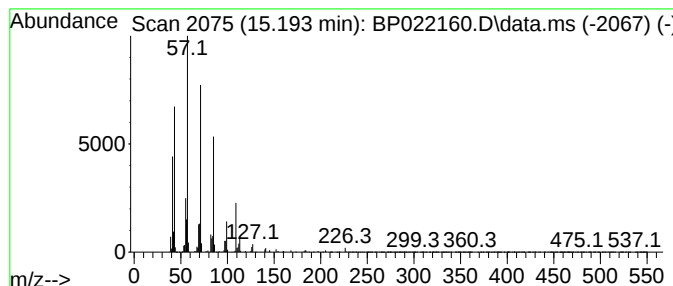
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BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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 4

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.193	64.49 ng/ul	7821210	Acenaphthene-d10	14.328	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexadecane	226	C16H34	000544-76-3	89
2	Hexadecane	226	C16H34	000544-76-3	89
3	Heptadecane	240	C17H36	000629-78-7	86
4	Dodecane, 4-methyl-	184	C13H28	006117-97-1	81
5	Tetradecane	198	C14H30	000629-59-4	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

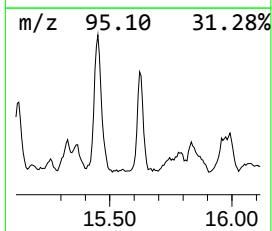
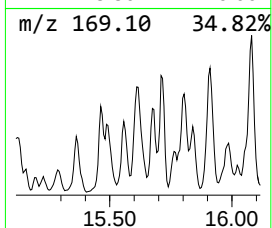
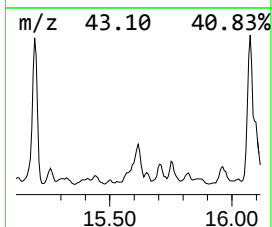
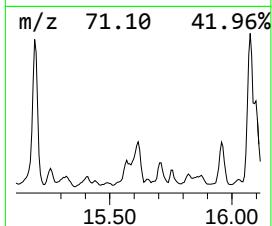
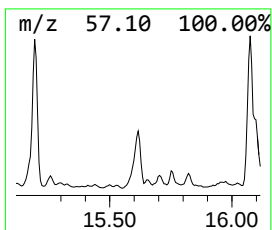
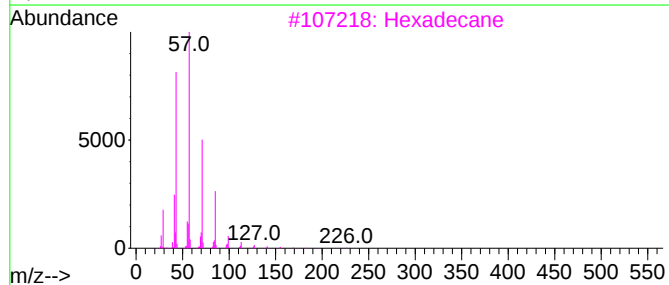
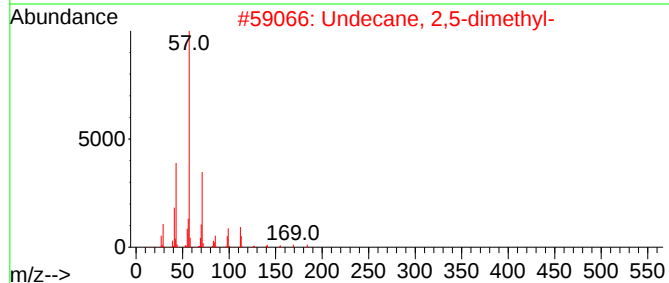
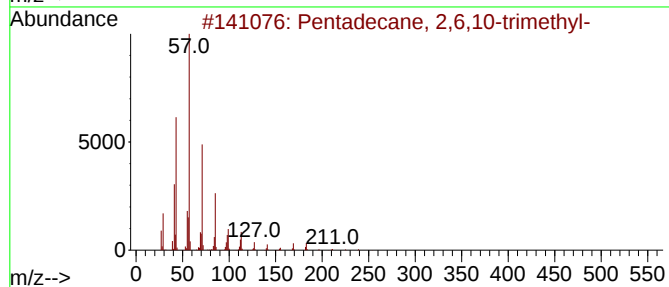
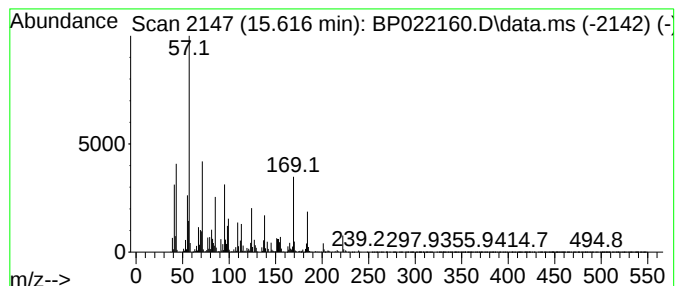
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ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 47 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
15.616	40.67 ng/ul	4933190	Acenaphthene-d10			14.328
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane, 2,6,10-trimethyl-		254	C18H38	003892-00-0	62
2	Undecane, 2,5-dimethyl-		184	C13H28	017301-22-3	46
3	Hexadecane		226	C16H34	000544-76-3	30
4	Dodecane		170	C12H26	000112-40-3	30
5	Dodecane		170	C12H26	000112-40-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

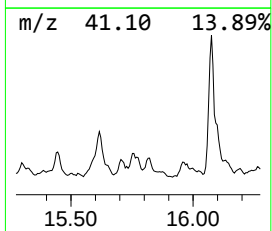
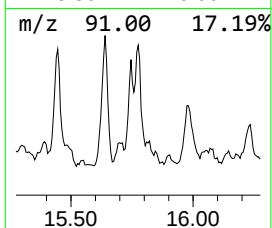
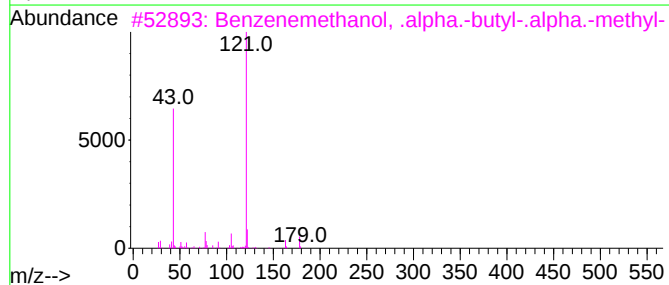
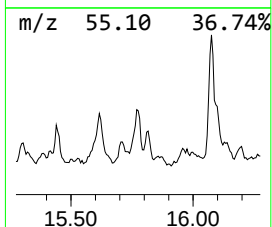
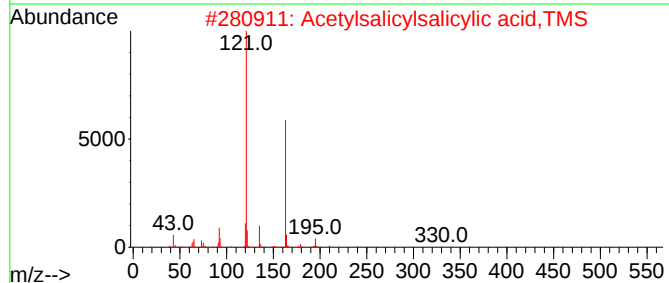
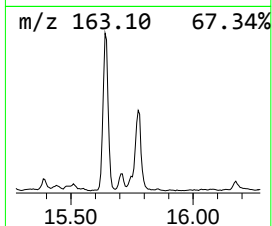
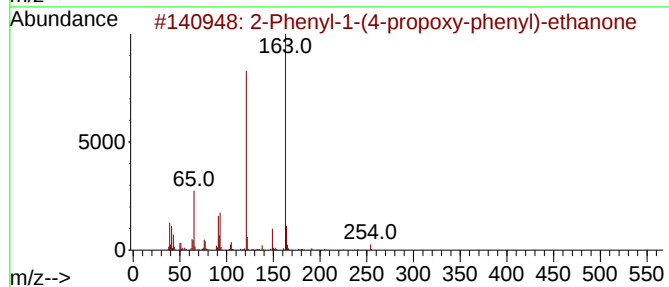
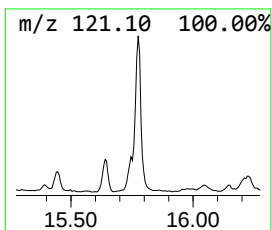
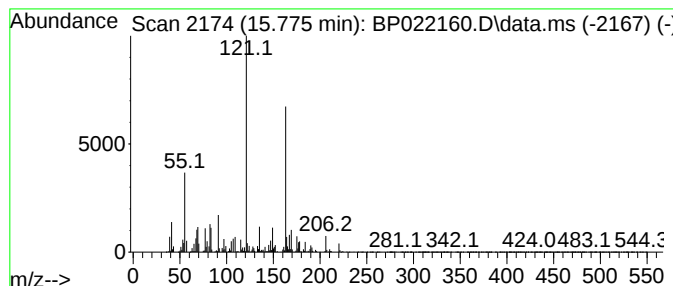
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Quant Title : SVOA CALIBRATION

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

Peak Number 48 2-Phenyl-1-(4-propoxy-pheny... Concentration Rank 49

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.775	13.49 ng/ul	1927190	Phenanthrene-d10	17.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Phenyl-1-(4-propoxy-phenyl)-et...	254	C17H18O2	063192-18-7	59
2			Acetylsalicylsalicylic acid,TMS	372	C19H20O6Si	1010472-34-3	59
3			Benzenemethanol, .alpha.-butyl-....	178	C12H18O	004396-98-9	38
4			Benzene, (1-methoxypropyl)-	150	C10H14O	059588-12-4	35
5			Benzene, (3-iodo-1-methoxybutyl)-	290	C11H15IO	1000327-38-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

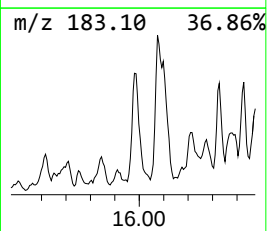
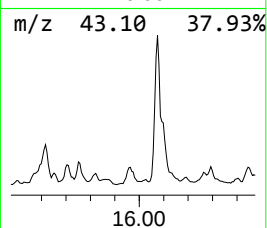
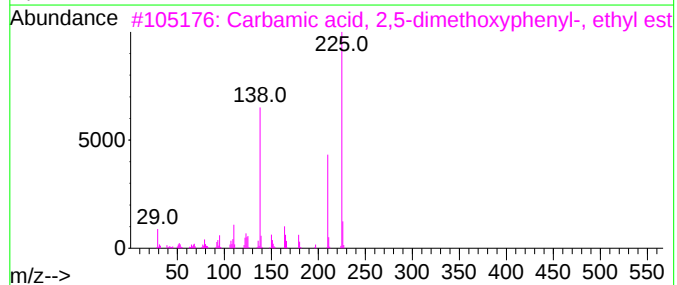
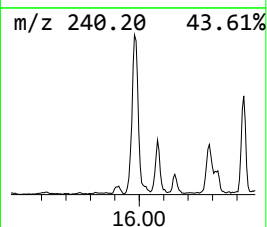
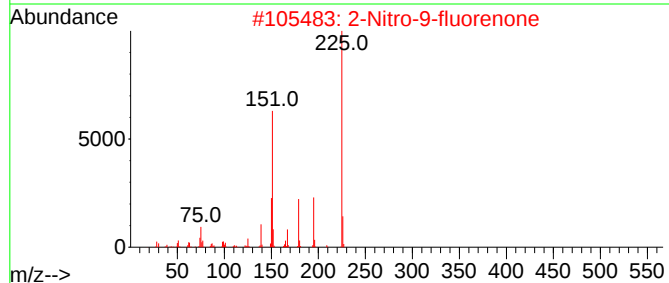
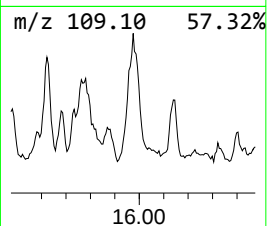
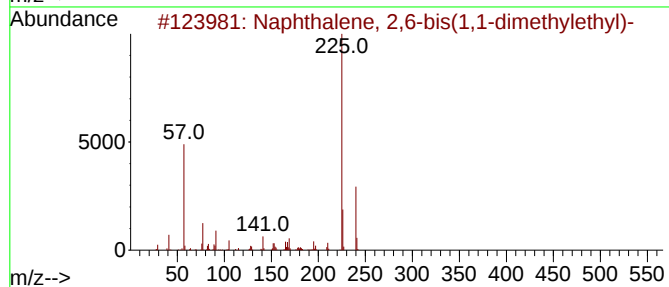
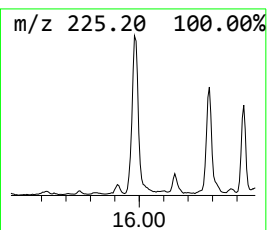
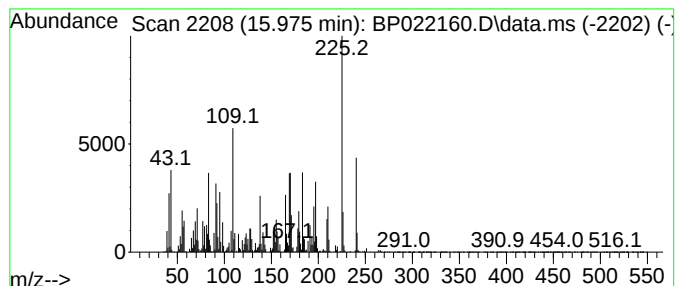
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Quant Title : SVOA CALIBRATION

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

Peak Number 49 unknown-06 Concentration Rank 30

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.975	19.69 ng/ul	2812920	Phenanthrene-d10	17.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,6-bis(1,1-dimethy...	240	C18H24	003905-64-4	38
2			2-Nitro-9-fluorenone	225	C13H7NO3	003096-52-4	30
3			Carbamic acid, 2,5-dimethoxyphen...	225	C11H15NO4	1000314-76-1	30
4			Acetamide, N-(2,5-dimethoxypheny...	225	C11H15NO4	1000307-26-1	30
5			(3-Isopropyl-5-methylamino-2,4,6...	225	C14H15N3	014204-00-3	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

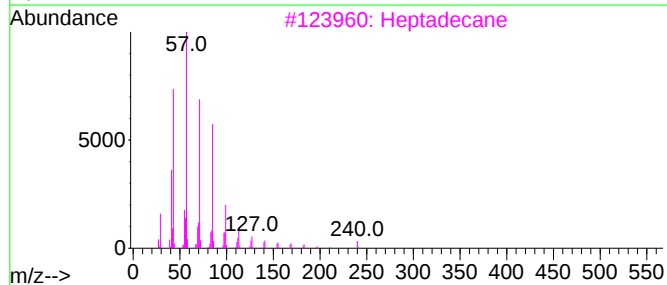
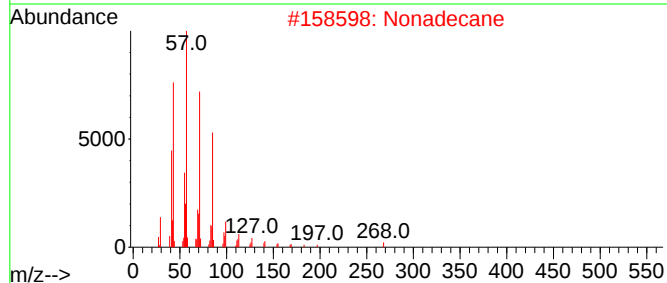
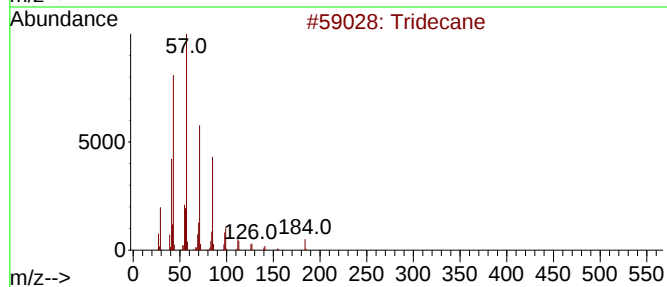
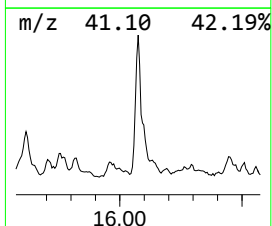
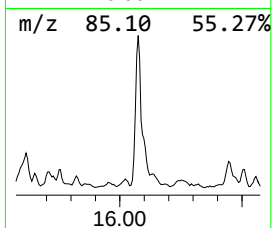
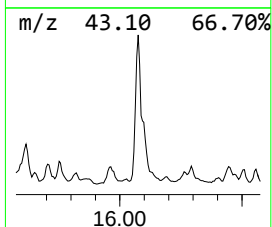
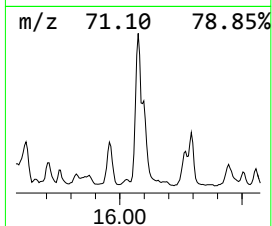
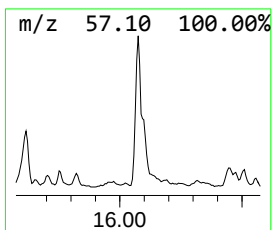
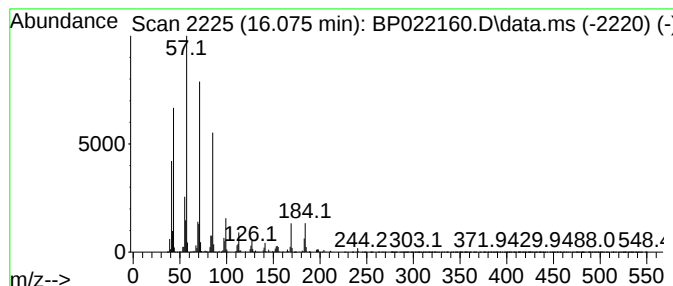
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 50 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.075	61.71 ng/ul	8814660	Phenanthrene-d10		17.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Tridecane		184	C13H28	000629-50-5	90
2	Nonadecane		268	C19H40	000629-92-5	87
3	Heptadecane		240	C17H36	000629-78-7	86
4	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	83
5	2-Bromo dodecane		248	C12H25Br	013187-99-0	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

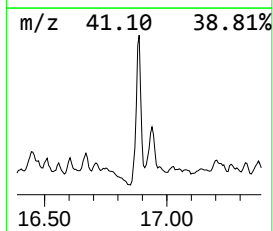
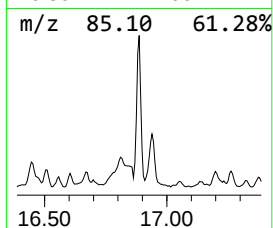
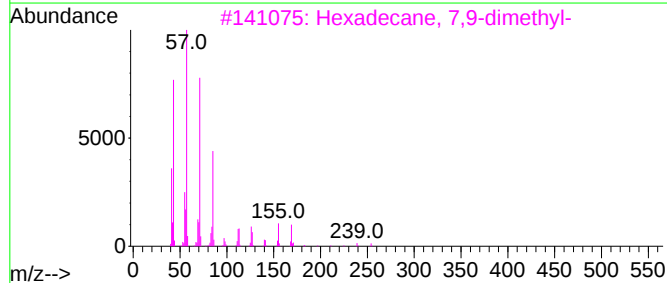
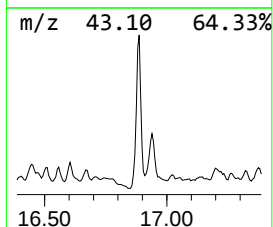
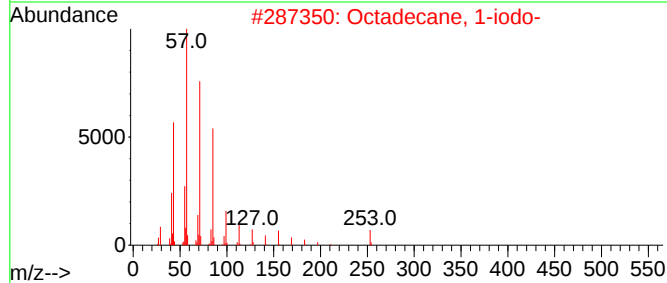
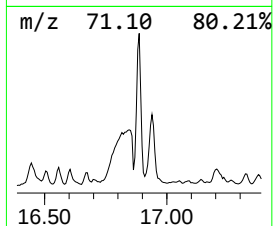
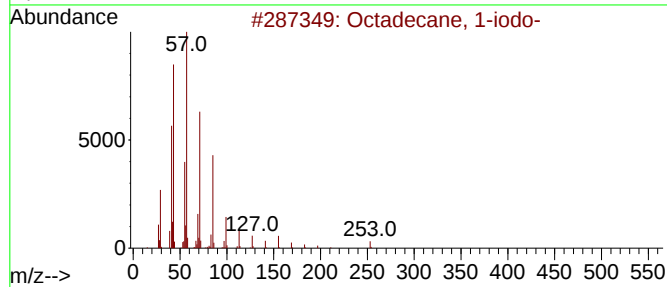
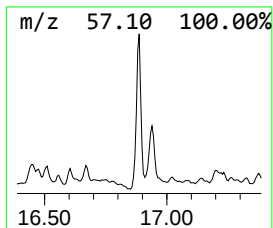
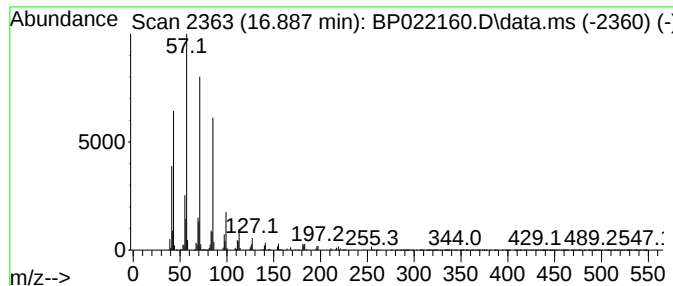
Instrument :
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ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 51 Octadecane, 1-iodo- Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD			R.T.
16.887	34.06 ng/ul	4865600	Phenanthrene-d10			17.145
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	86
2	Octadecane, 1-iodo-		380	C18H37I	000629-93-6	86
3	Hexadecane, 7,9-dimethyl-		254	C18H38	021164-95-4	76
4	Eicosane		282	C20H42	000112-95-8	72
5	Heptadecane		240	C17H36	000629-78-7	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

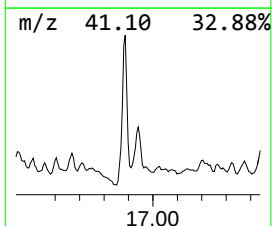
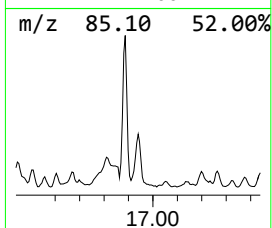
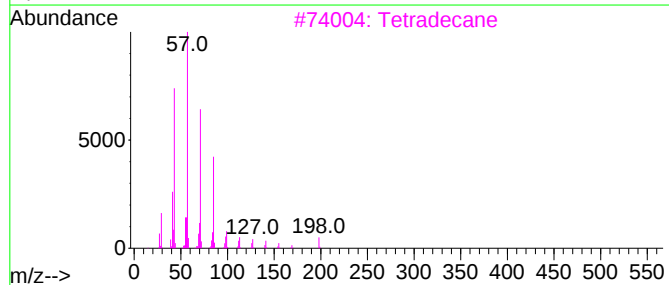
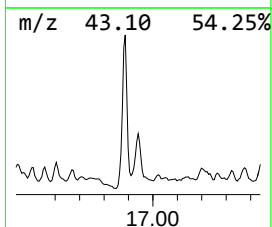
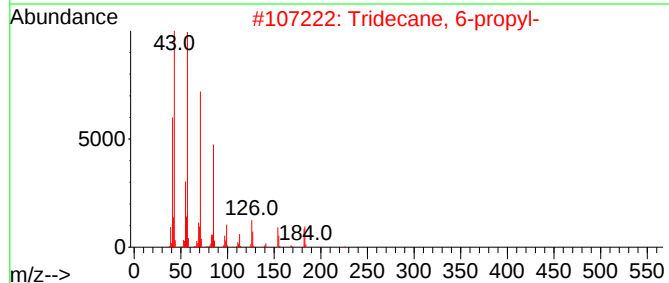
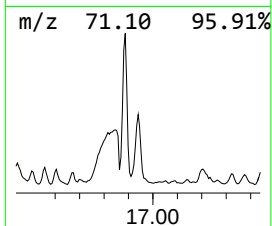
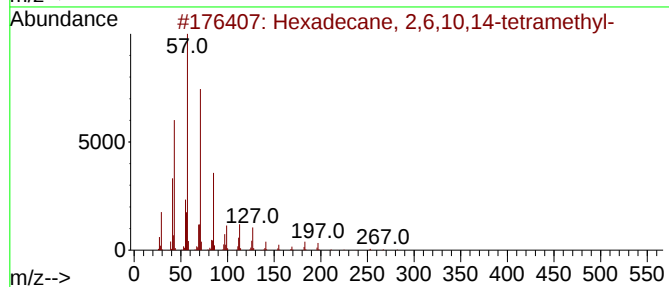
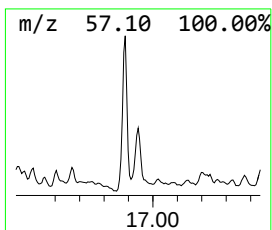
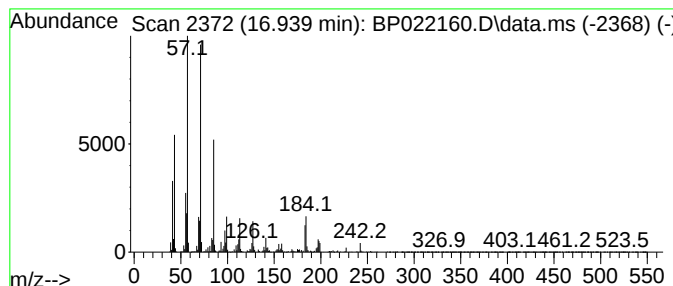
Instrument :
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ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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 41

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.939	16.48 ng/ul	2354640	Phenanthrene-d10		17.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane, 2,6,10,14-tetramethyl-		282	C20H42	000638-36-8	96
2	Tridecane, 6-propyl-		226	C16H34	055045-10-8	87
3	Tetradecane		198	C14H30	000629-59-4	81
4	Dodecane		170	C12H26	000112-40-3	81
5	Heptadecane, 2,6,10,15-tetramethyl-		296	C21H44	054833-48-6	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

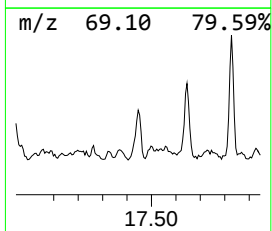
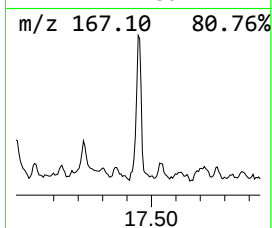
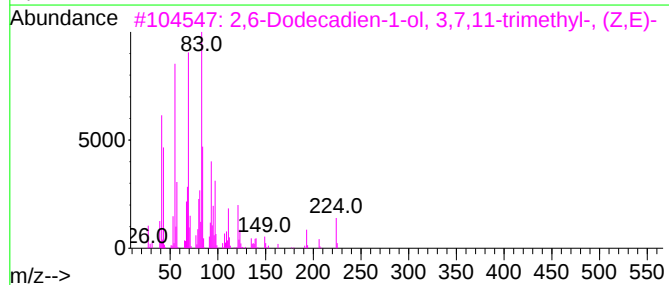
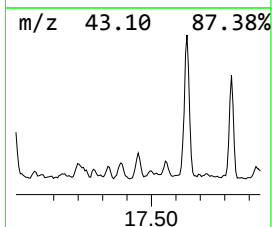
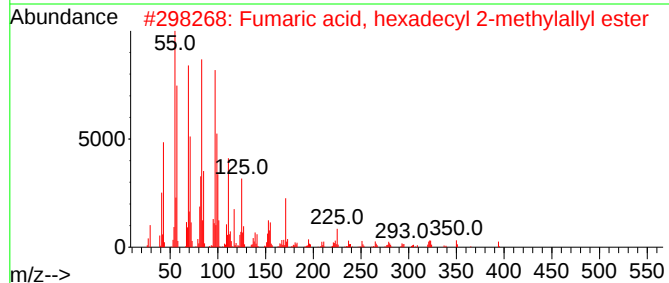
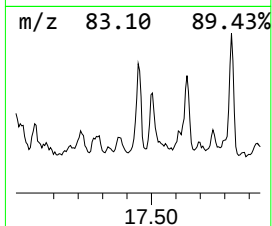
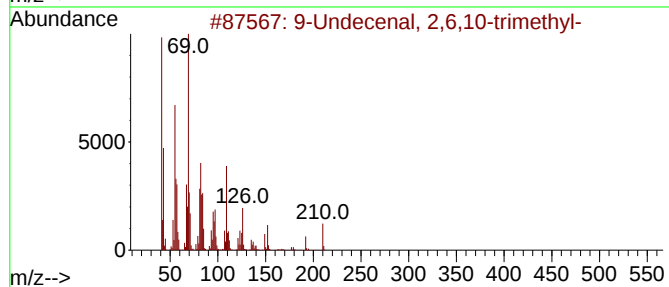
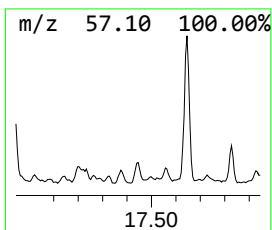
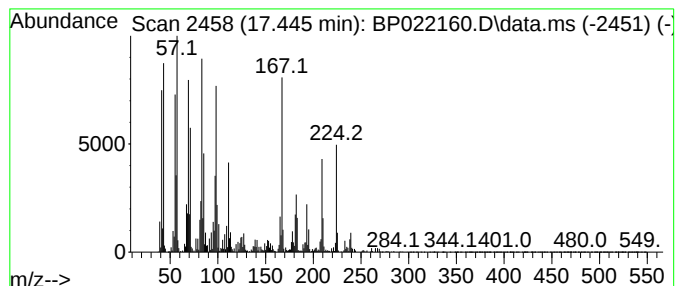
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Quant Title : SVOA CALIBRATION

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

Peak Number 53 unknown-07 Concentration Rank 31

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.445	19.26 ng/ul	2751800	Phenanthrene-d10	17.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Undecenal, 2,6,10-trimethyl-	210	C14H26O	000141-13-9	46		
2	Fumaric acid, hexadecyl 2-methyl...	394	C24H42O4	1000330-55-5	42		
3	2,6-Dodecadien-1-ol, 3,7,11-trim...	224	C15H28O	020576-58-3	41		
4	3-Pentenal, 4-methyl-	98	C6H10O	005362-50-5	35		
5	Cyclopropane, 1-methyl-1-(2-meth...	238	C17H34	041977-41-7	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

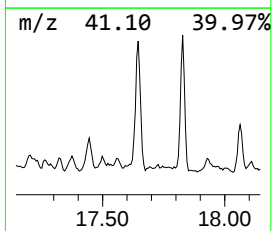
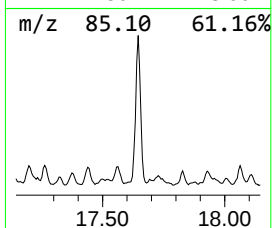
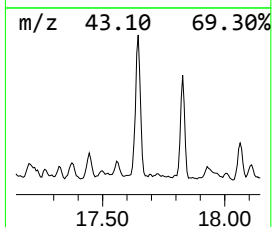
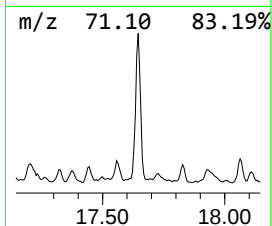
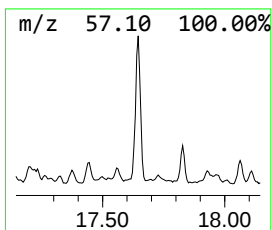
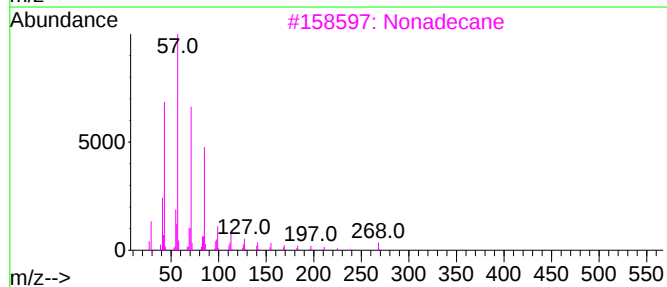
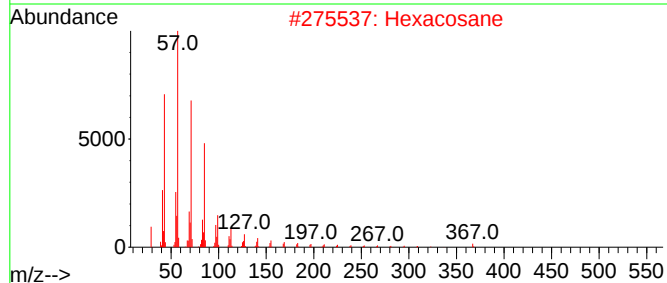
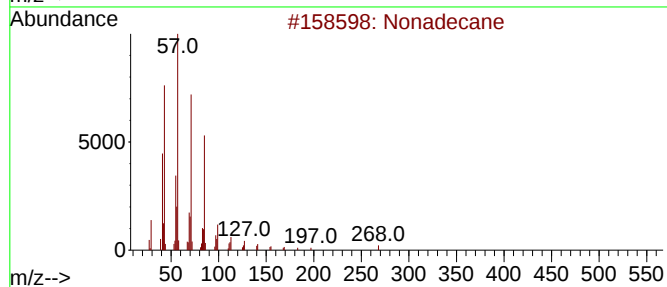
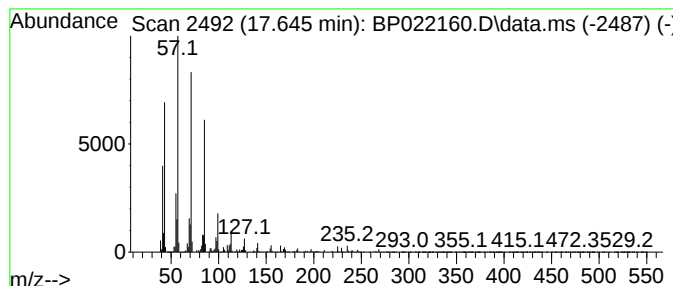
Instrument :
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ClientSampleId :
DCVW3

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Quant Title : SVOA CALIBRATION

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

Peak Number 54 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.645	38.21 ng/ul	5458600	Phenanthrene-d10		17.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	96
2	Hexacosane		366	C26H54	000630-01-3	91
3	Nonadecane		268	C19H40	000629-92-5	91
4	Tetradecane		198	C14H30	000629-59-4	86
5	Tridecane, 6-propyl-		226	C16H34	055045-10-8	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

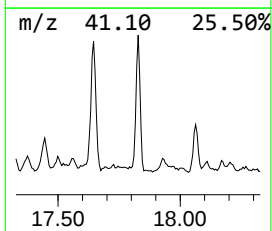
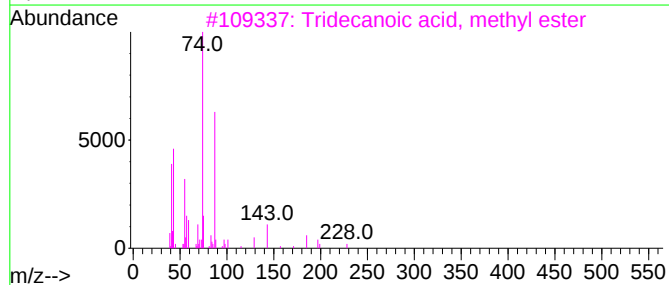
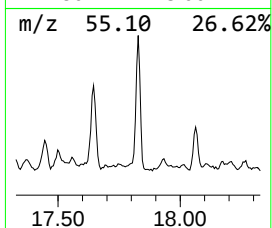
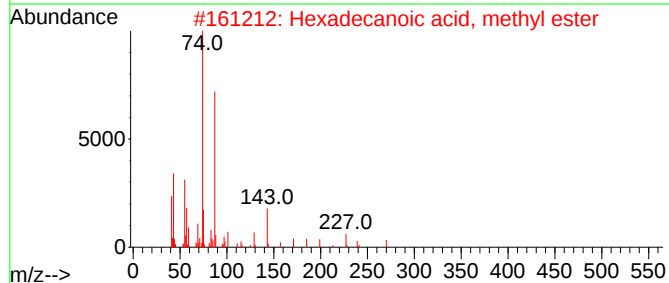
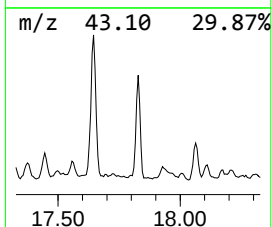
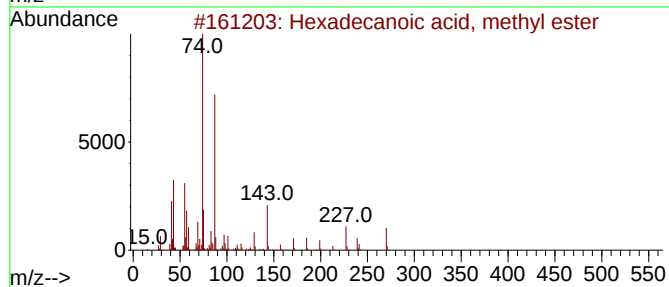
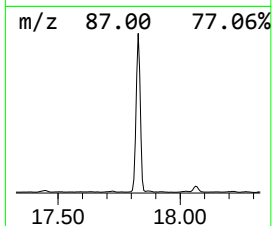
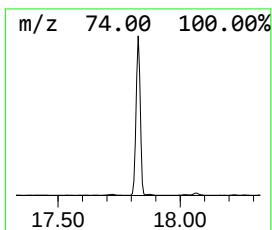
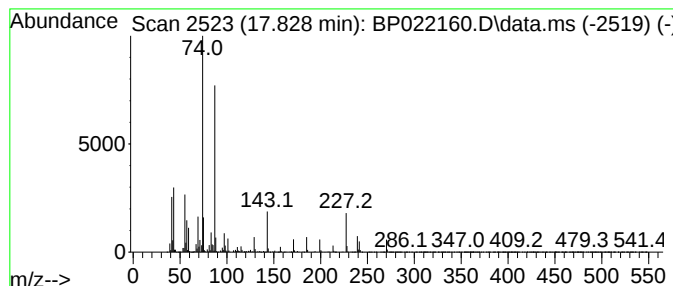
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 55 Hexadecanoic acid, methyl e... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.828	42.21 ng/ul	6029760	Phenanthrene-d10		17.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecanoic acid, methyl ester		270	C17H34O2	000112-39-0	97
2	Hexadecanoic acid, methyl ester		270	C17H34O2	000112-39-0	96
3	Tridecanoic acid, methyl ester		228	C14H28O2	001731-88-0	93
4	Hexadecanoic acid, methyl ester		270	C17H34O2	000112-39-0	90
5	Methyl tetradecanoate		242	C15H30O2	000124-10-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

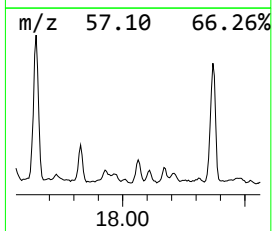
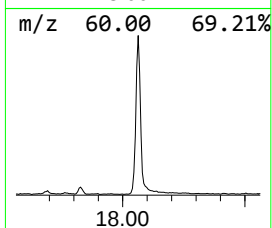
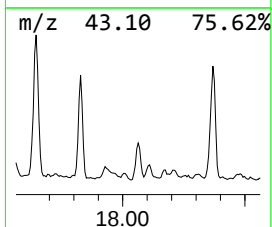
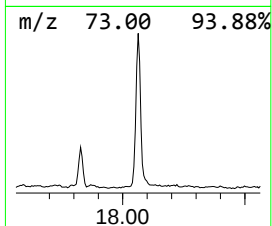
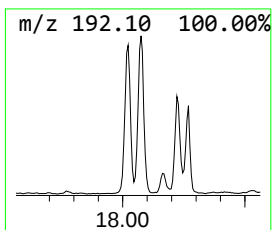
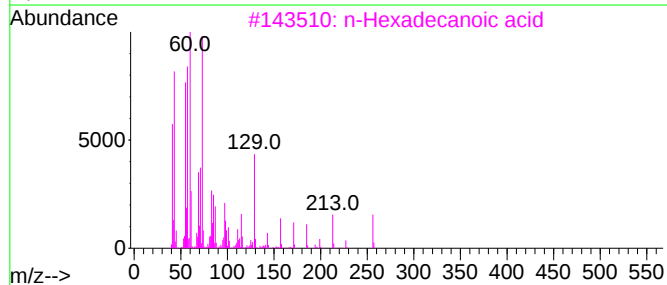
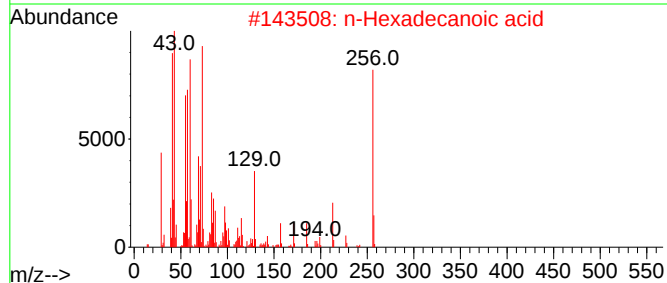
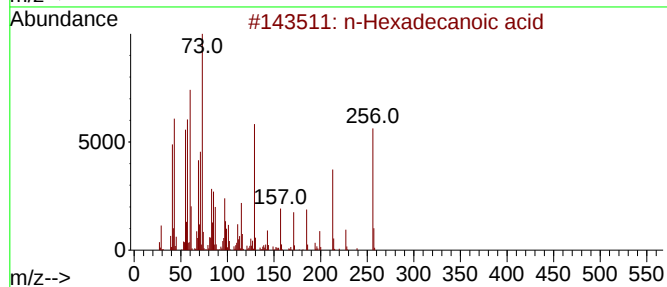
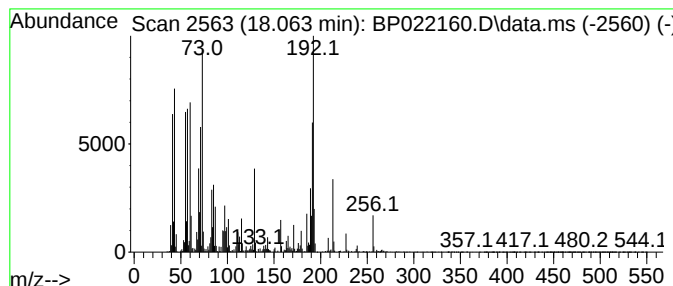
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 56 n-Hexadecanoic acid Concentration Rank 35

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.063	17.59 ng/ul	2512090	Phenanthrene-d10	17.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	86
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	84
4	Phenanthrene, 2-methyl-	192	C15H12	002531-84-2	55
5	Anthracene, 1-methyl-	192	C15H12	000610-48-0	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

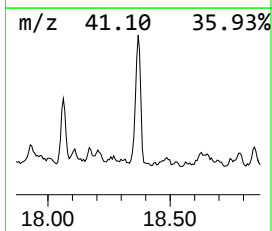
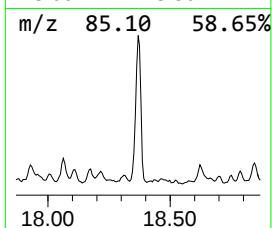
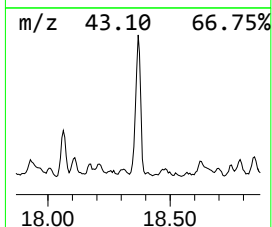
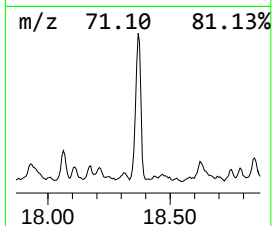
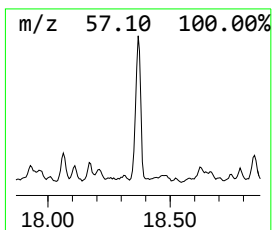
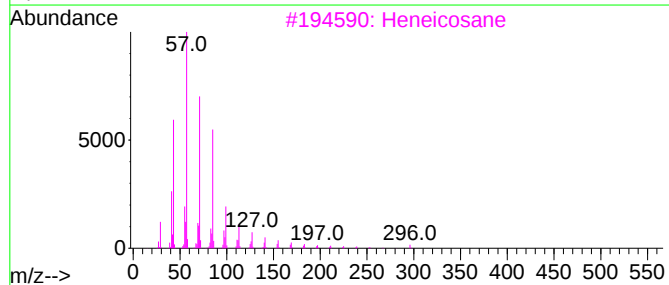
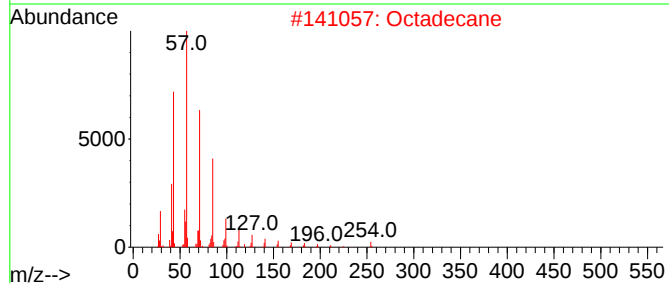
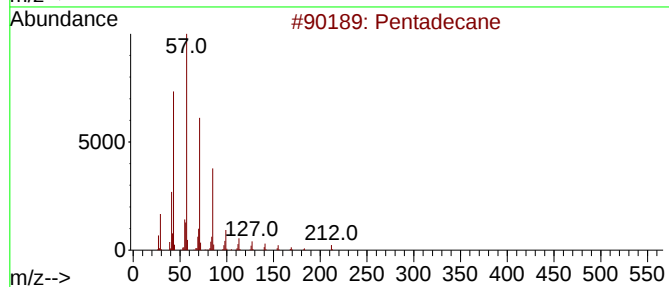
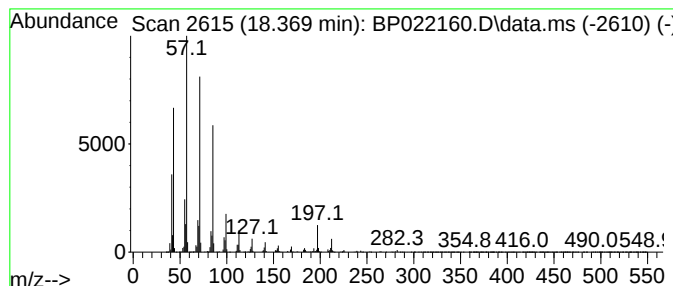
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 57 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.369	28.04 ng/ul	4004940	Phenanthrene-d10		17.145	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Pentadecane		212	C15H32	000629-62-9	90
2	Octadecane		254	C18H38	000593-45-3	87
3	Heneicosane		296	C21H44	000629-94-7	87
4	Heneicosane		296	C21H44	000629-94-7	87
5	Pentadecane		212	C15H32	000629-62-9	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

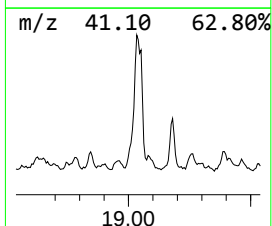
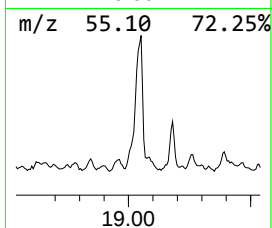
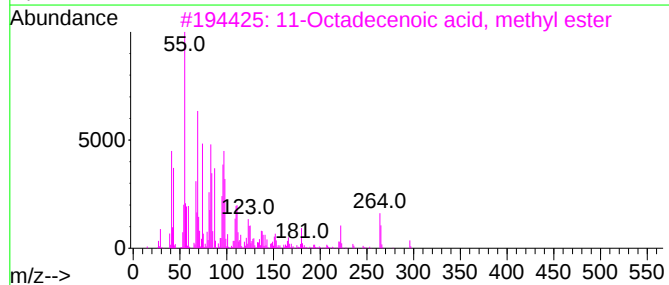
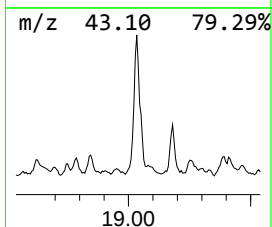
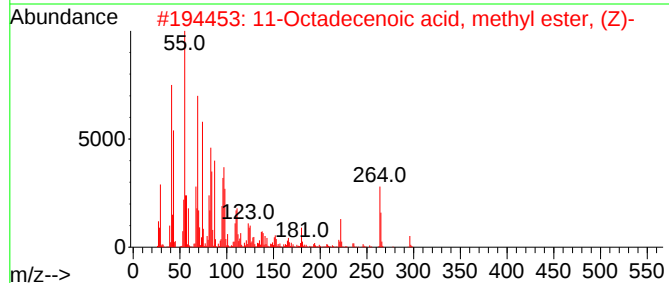
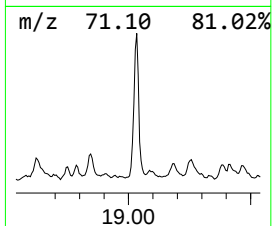
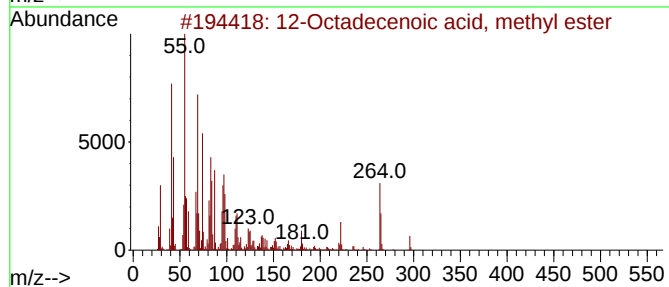
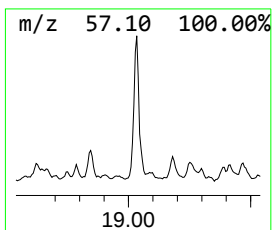
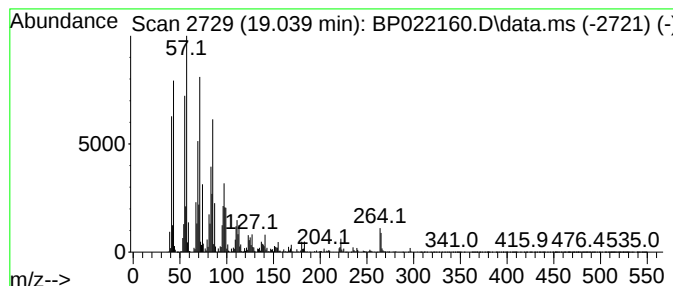
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Quant Title : SVOA CALIBRATION

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

Peak Number 58 12-Octadecenoic acid, methyl... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.039	56.90 ng/ul	8127660	Phenanthrene-d10	17.145

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			12-Octadecenoic acid, methyl ester	296	C19H36O2	056554-46-2	96
2			11-Octadecenoic acid, methyl est...	296	C19H36O2	001937-63-9	96
3			11-Octadecenoic acid, methyl ester	296	C19H36O2	052380-33-3	96
4			9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	95
5			9-Octadecenoic acid, methyl este...	296	C19H36O2	001937-62-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

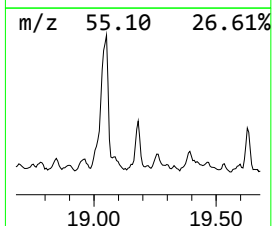
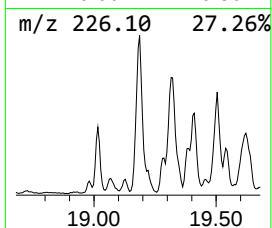
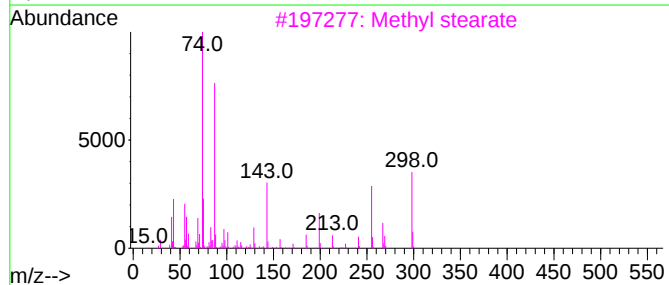
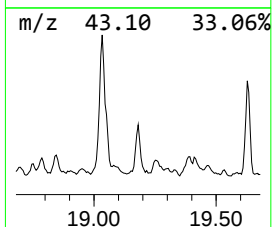
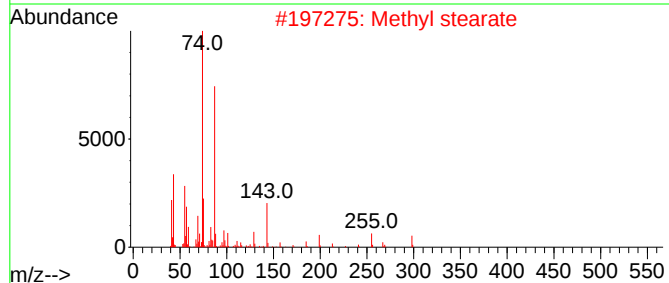
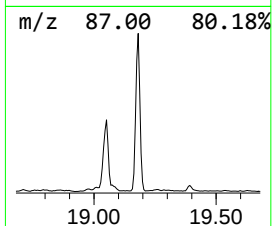
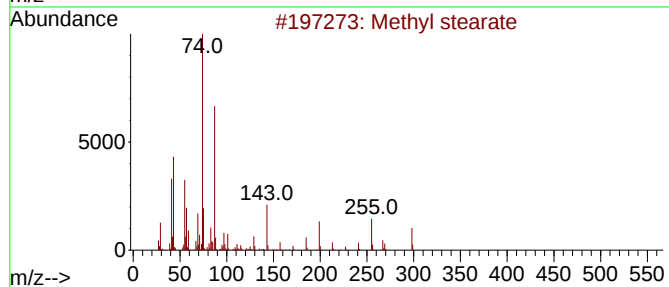
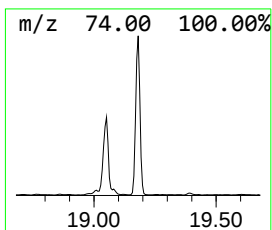
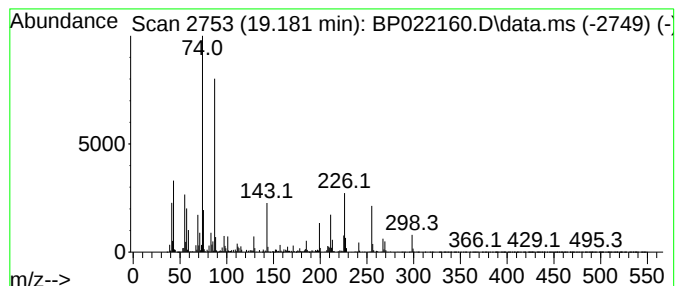
Instrument :
BNA_P
ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 59 Methyl stearate Concentration Rank 34

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.180	18.65 ng/ul	2663370	Phenanthrene-d10	17.145	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl stearate	298	C19H38O2	000112-61-8	99
2	Methyl stearate	298	C19H38O2	000112-61-8	97
3	Methyl stearate	298	C19H38O2	000112-61-8	96
4	Methyl stearate	298	C19H38O2	000112-61-8	95
5	Methyl stearate	298	C19H38O2	000112-61-8	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

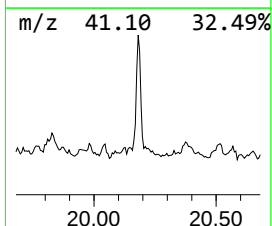
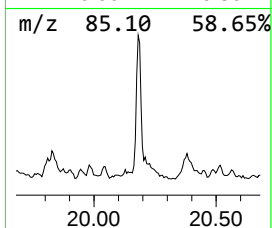
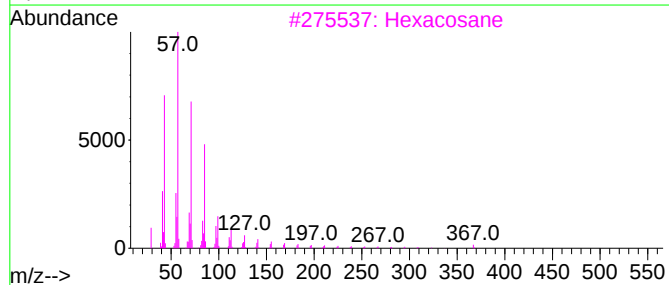
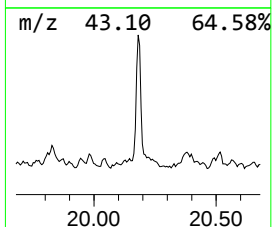
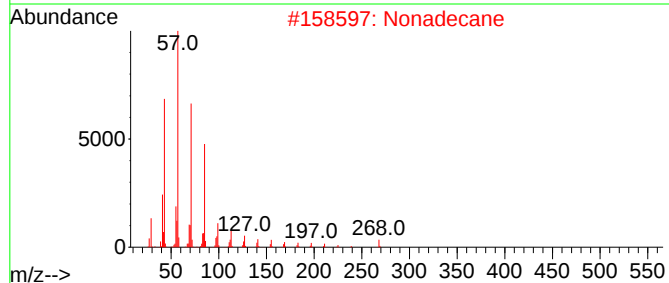
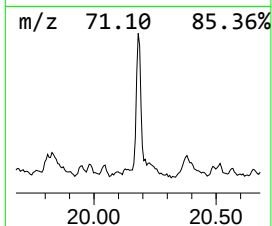
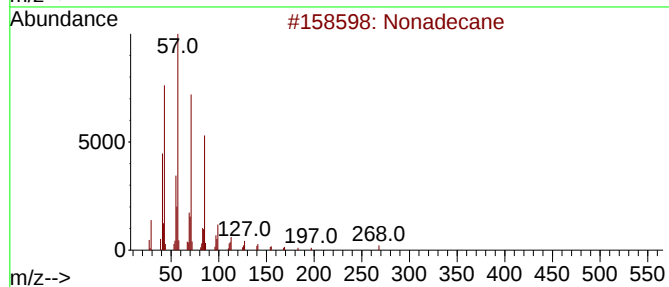
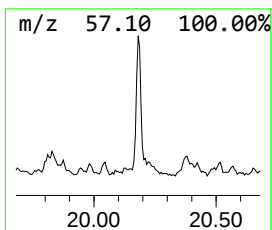
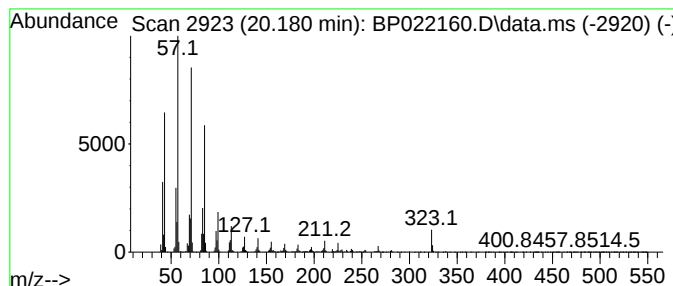
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Quant Title : SVOA CALIBRATION

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

Peak Number 60 (DEL) Alkane: Straight-Chai... Concentration Rank 29

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.180	19.87 ng/ul	2462360	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Nonadecane	268	C19H40	000629-92-5	90
3			Hexacosane	366	C26H54	000630-01-3	87
4			Tricosane, 2-methyl-	338	C24H50	001928-30-9	87
5			Heptadecane, 9-octyl-	352	C25H52	007225-64-1	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCVW3

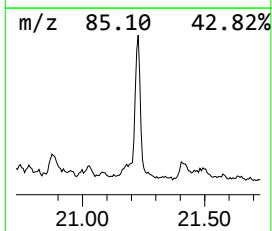
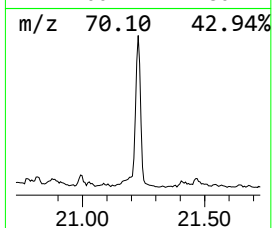
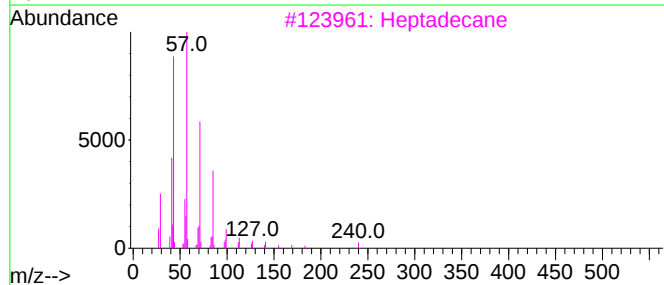
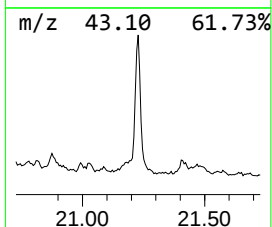
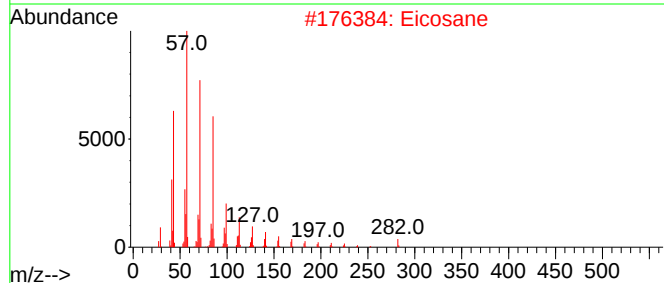
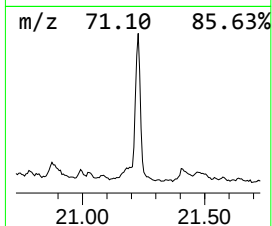
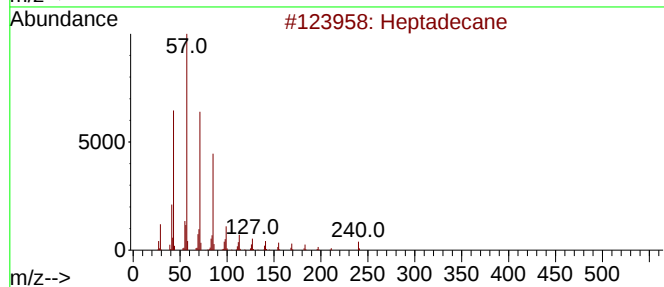
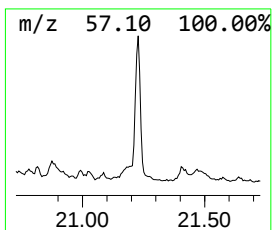
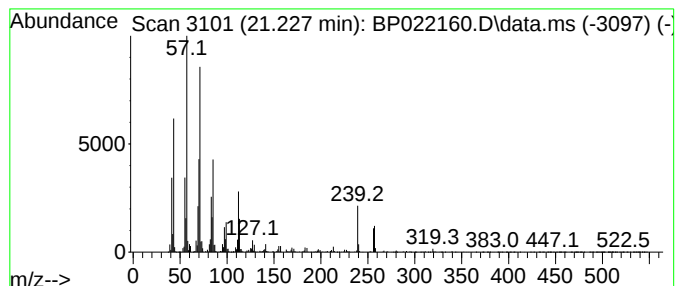
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Quant Title : SVOA CALIBRATION

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

Peak Number 61 (DEL) Alkane: Straight-Chai... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.227	34.51 ng/ul	4277520	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane	240	C17H36	000629-78-7	80
2			Eicosane	282	C20H42	000112-95-8	70
3			Heptadecane	240	C17H36	000629-78-7	60
4			Pentadecane	212	C15H32	000629-62-9	60
5			Heptadecane, 7-methyl-	254	C18H38	020959-33-5	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

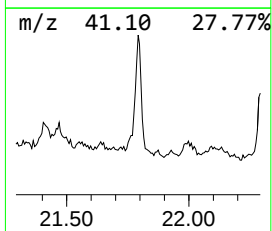
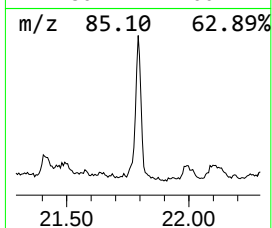
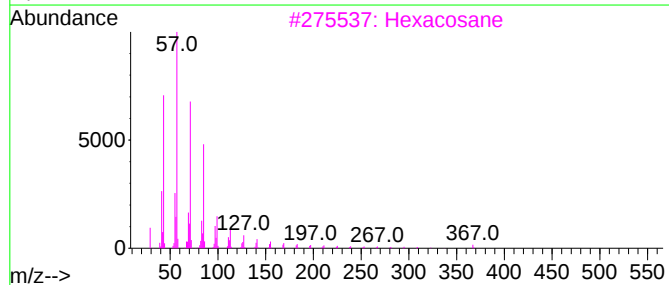
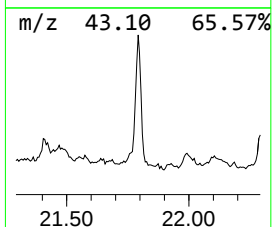
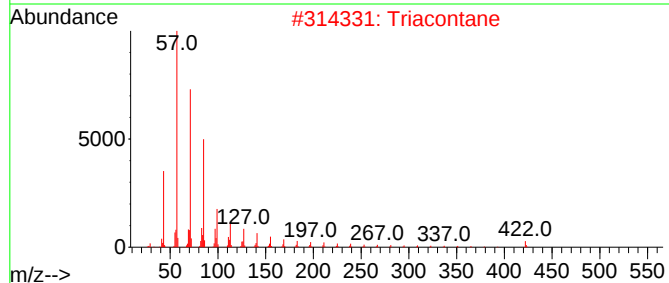
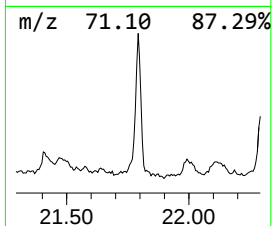
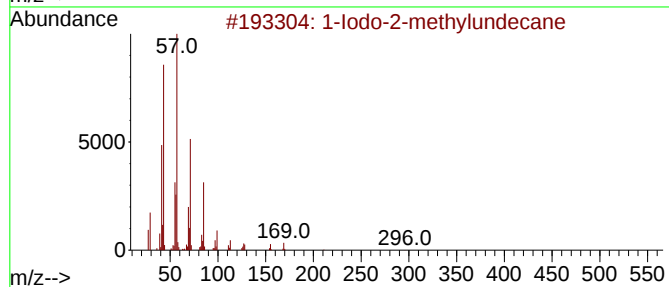
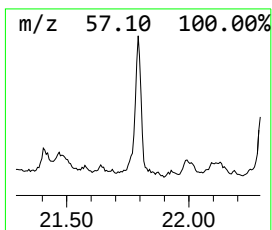
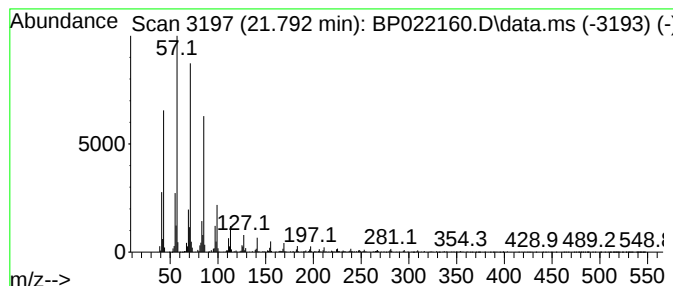
Instrument :
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ClientSampleId :
DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 62 1-Iodo-2-methylundecane Concentration Rank 42

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.792	15.93 ng/ul	1974000	Chrysene-d12	21.545	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Iodo-2-methylundecane	296	C12H25I	073105-67-6	91
2	Triacontane	422	C30H62	000638-68-6	91
3	Hexacosane	366	C26H54	000630-01-3	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Hexadecane, 3-methyl-	240	C17H36	006418-43-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
Data File : BP022160.D
Acq On : 02 Oct 2024 13:25
Operator : RC/JU
Sample : P4016-12
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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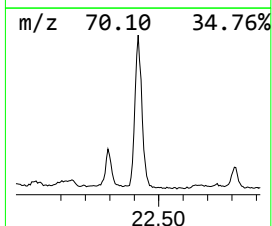
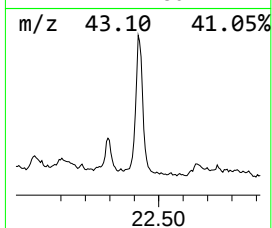
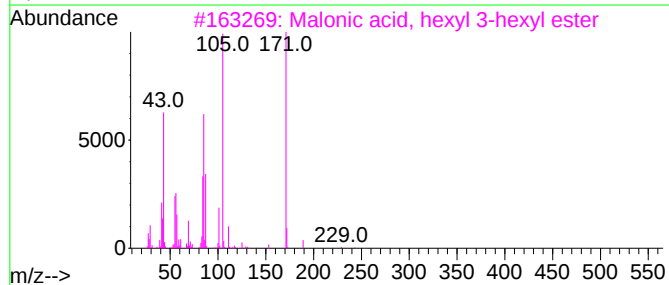
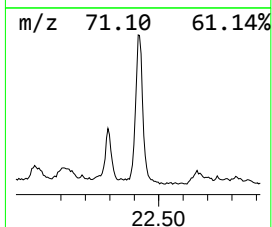
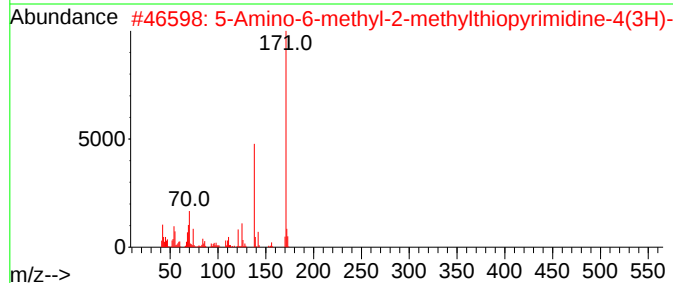
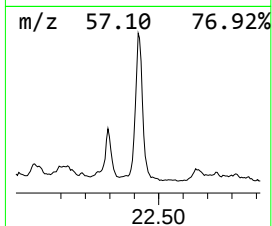
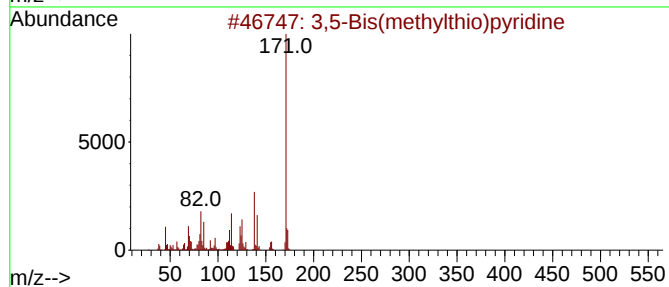
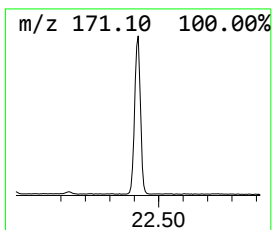
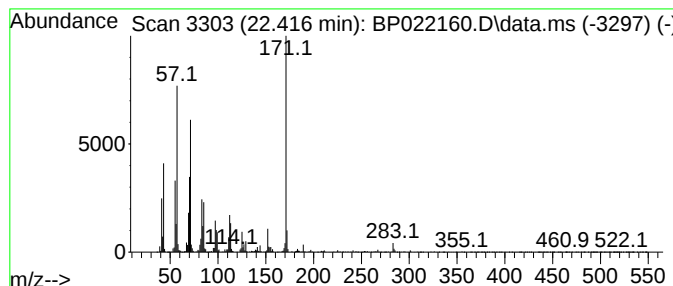
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Quant Title : SVOA CALIBRATION

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

Peak Number 63 unknown-08 Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.416	36.70 ng/ul	4548960	Chrysene-d12	21.545

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			3,5-Bis(methylthio)pyridine	171	C7H9NS2	070999-08-5	47
2			5-Amino-6-methyl-2-methylthiopyr...	171	C6H9N3OS	342402-52-2	38
3			Malonic acid, hexyl 3-hexyl ester	272	C15H28O4	1000349-12-2	38
4			Nonanedioic acid, bis(2-ethylhex...	412	C25H48O4	000103-24-2	38
5			Succinic acid, di(3-oxobut-2-yl)...	258	C12H18O6	1000349-59-0	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100224\
 Data File : BP022160.D
 Acq On : 02 Oct 2024 13:25
 Operator : RC/JU
 Sample : P4016-12
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCVW3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092424.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
(DEL) Alkane: S...	5.752	14.0	ng/ul	3022490	1	7.699	4307400	20.0
(DEL) Alkane: S...	6.711	17.5	ng/ul	3760300	1	7.699	4307400	20.0
(DEL) Alkane: S...	6.769	11.7	ng/ul	2521220	1	7.699	4307400	20.0
Benzene, 1-ethy...	6.811	19.0	ng/ul	4094820	1	7.699	4307400	20.0
(DEL) Alkane: S...	7.346	89.5	ng/ul	19275700	1	7.699	4307400	20.0
(DEL) Alkane: S...	7.622	12.4	ng/ul	2668680	1	7.699	4307400	20.0
(DEL) Alkane: S...	7.769	14.1	ng/ul	3032510	1	7.699	4307400	20.0
Mesitylene	7.816	16.8	ng/ul	3627030	1	7.699	4307400	20.0
(DEL) Alkane: C...	7.987	9.7	ng/ul	2094290	1	7.699	4307400	20.0
Cyclohexanone, ...	8.128	13.8	ng/ul	2979600	1	7.699	4307400	20.0
Benzene, 1,4-di...	8.181	14.3	ng/ul	3084590	1	7.699	4307400	20.0
Benzene, 1-meth...	8.234	24.9	ng/ul	5374460	1	7.699	4307400	20.0
(DEL) Alkane: S...	8.287	10.7	ng/ul	2312000	1	7.699	4307400	20.0
Benzene, 1-ethy...	8.334	45.6	ng/ul	9824430	1	7.699	4307400	20.0
(DEL) Alkane: S...	8.463	16.8	ng/ul	3627610	1	7.699	4307400	20.0
o-Cymene	8.693	18.7	ng/ul	4033710	1	7.699	4307400	20.0
(DEL) Alkane: S...	8.922	79.9	ng/ul	17211000	1	7.699	4307400	20.0
unknown-01	9.040	27.1	ng/ul	5825670	1	7.699	4307400	20.0
(DEL) Alkane: S...	10.975	4.8	ng/ul	6954050	2	10.463	28665200	20.0
(DEL) Alkane: S...	11.728	3.2	ng/ul	4612000	2	10.463	28665200	20.0
unknown-02	12.563	36.8	ng/ul	4463230	3	14.328	2425730	20.0
unknown-03	12.646	17.0	ng/ul	2067220	3	14.328	2425730	20.0
1H-Pyrazole, 4,...	12.887	17.0	ng/ul	2063460	3	14.328	2425730	20.0
unknown-04	13.369	47.8	ng/ul	5798210	3	14.328	2425730	20.0
Naphthalene, 2,...	13.499	54.2	ng/ul	6573840	3	14.328	2425730	20.0
Naphthalene, 1,...	13.646	35.9	ng/ul	4349970	3	14.328	2425730	20.0
(DEL) Alkane: S...	13.804	25.5	ng/ul	3092450	3	14.328	2425730	20.0
Naphthalene, 1,...	13.869	21.4	ng/ul	2601780	3	14.328	2425730	20.0
unknown-05	14.128	20.2	ng/ul	2453650	3	14.328	2425730	20.0
(DEL) Alkane: S...	14.240	70.5	ng/ul	8548470	3	14.328	2425730	20.0
9-Methyl-S-octa...	14.463	21.9	ng/ul	2654400	3	14.328	2425730	20.0
Naphthalene, 1,...	14.734	22.9	ng/ul	2782520	3	14.328	2425730	20.0
Naphthalene, 2,...	14.798	21.2	ng/ul	2572510	3	14.328	2425730	20.0
(DEL) Alkane: S...	14.857	22.6	ng/ul	2746070	3	14.328	2425730	20.0
Naphthalene, 1,...	15.098	30.3	ng/ul	3677840	3	14.328	2425730	20.0
(DEL) Alkane: S...	15.193	64.5	ng/ul	7821210	3	14.328	2425730	20.0
(DEL) Alkane: S...	15.616	40.7	ng/ul	4933190	3	14.328	2425730	20.0
2-Phenyl-1-(4-p...	15.775	13.5	ng/ul	1927190	4	17.145	2856890	20.0
unknown-06	15.975	19.7	ng/ul	2812920	4	17.145	2856890	20.0
(DEL) Alkane: S...	16.075	61.7	ng/ul	8814660	4	17.145	2856890	20.0
Octadecane, 1-i...	16.887	34.1	ng/ul	4865600	4	17.145	2856890	20.0
(DEL) Alkane: S...	16.939	16.5	ng/ul	2354640	4	17.145	2856890	20.0
unknown-07	17.445	19.3	ng/ul	2751800	4	17.145	2856890	20.0
(DEL) Alkane: S...	17.645	38.2	ng/ul	5458600	4	17.145	2856890	20.0
Hexadecanoic ac...	17.828	42.2	ng/ul	6029760	4	17.145	2856890	20.0
n-Hexadecanoic ...	18.063	17.6	ng/ul	2512090	4	17.145	2856890	20.0
(DEL) Alkane: S...	18.369	28.0	ng/ul	4004940	4	17.145	2856890	20.0
12-Octadecenoic...	19.039	56.9	ng/ul	8127660	4	17.145	2856890	20.0
Methyl stearate	19.180	18.6	ng/ul	2663370	4	17.145	2856890	20.0
(DEL) Alkane: S...	20.180	19.9	ng/ul	2462360	5	21.545	2478770	20.0
(DEL) Alkane: S...	21.227	34.5	ng/ul	4277520	5	21.545	2478770	20.0
1-Iodo-2-methyl...	21.792	15.9	ng/ul	1974000	5	21.545	2478770	20.0

Instrument :

BNA_P

ClientSampleId :

DCVW3