

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
 Data File : BM047748.D
 Acq On : 01 Oct 2024 18:21
 Operator : RC/JU
 Sample : P4018-08DL 30X
 Misc :
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 DDCB4DL

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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047748.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.005	168	173	181	rVB	60943	87354	0.60%	0.163%
2	5.834	478	484	491	rVB	58996	88398	0.61%	0.165%
3	6.810	646	650	655	rVB2	50422	78539	0.54%	0.146%
4	6.869	655	660	663	rBV	62294	92275	0.63%	0.172%
5	6.975	671	678	684	rVV2	109918	204385	1.40%	0.381%
6	7.040	684	689	698	rVB	59943	96547	0.66%	0.180%
7	7.240	720	723	727	rVV	61920	90820	0.62%	0.169%
8	7.446	752	758	767	rVV2	347270	635597	4.36%	1.185%
9	7.810	812	820	826	rBV	819891	1353305	9.28%	2.522%
10	7.881	826	832	837	rBV4	82650	145885	1.00%	0.272%
11	7.928	837	840	848	rVB2	50663	84516	0.58%	0.158%
12	8.110	867	871	878	rVB3	44436	84597	0.58%	0.158%
13	8.351	907	912	918	rVB2	91269	163555	1.12%	0.305%
14	8.410	918	922	925	rVB	68420	85677	0.59%	0.160%
15	8.452	925	929	931	rBV2	82363	116149	0.80%	0.216%
16	8.475	931	933	937	rVB	89810	111365	0.76%	0.208%
17	8.581	945	951	955	rBV	116025	194165	1.33%	0.362%
18	8.616	955	957	965	rVB2	59356	97525	0.67%	0.182%
19	8.763	978	982	986	rBV	57874	89600	0.61%	0.167%
20	8.816	986	991	998	rVB2	72411	119329	0.82%	0.222%
21	8.922	999	1009	1013	rBV3	89194	199492	1.37%	0.372%
22	9.046	1020	1030	1035	rBV	620348	919055	6.30%	1.713%
23	9.169	1041	1051	1056	rBV9	72942	203812	1.40%	0.380%
24	9.251	1056	1065	1068	rBV2	44723	94720	0.65%	0.177%
25	9.304	1068	1074	1082	rVB5	41763	130411	0.89%	0.243%
26	9.604	1118	1125	1132	rVB4	31511	81476	0.56%	0.152%
27	9.698	1132	1141	1145	rBV4	50507	114050	0.78%	0.213%
28	9.746	1145	1149	1154	rVV4	61333	114048	0.78%	0.213%
29	9.798	1154	1158	1166	rVB4	36432	97122	0.67%	0.181%
30	9.893	1166	1174	1180	rVB	85413	187923	1.29%	0.350%
31	9.963	1180	1186	1190	rBV2	69938	116314	0.80%	0.217%
32	10.040	1196	1199	1203	rVV	66625	108046	0.74%	0.201%
33	10.146	1212	1217	1222	rBV2	56201	102525	0.70%	0.191%
34	10.304	1238	1244	1249	rBV2	64725	109930	0.75%	0.205%
35	10.510	1275	1279	1283	rVV2	43003	82094	0.56%	0.153%
36	10.598	1284	1294	1301	rVV4	1640087	3203402	21.96%	5.970%

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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_M\Methods\SFAM-EPA-BM092724.MA.M
 Title : SVOA CALIBRATION

37	10.651	1301	1303	1310	rVV2	71686	133779	0.92%	0.249%
38	10.728	1310	1316	1322	rVV4	249181	521406	3.57%	0.972%
39	10.775	1322	1324	1329	rVV	86346	130382	0.89%	0.243%
40	10.987	1354	1360	1368	rBV	267885	454061	3.11%	0.846%

41	11.116	1376	1382	1388	rVB	68591	132557	0.91%	0.247%
42	11.298	1405	1413	1418	rBV3	75333	187107	1.28%	0.349%
43	11.492	1442	1446	1449	rVV3	47966	97646	0.67%	0.182%
44	11.528	1449	1452	1461	rVB5	62540	124875	0.86%	0.233%
45	11.640	1465	1471	1475	rBV	153731	273988	1.88%	0.511%

46	11.704	1475	1482	1490	rVB2	286226	566992	3.89%	1.057%
47	11.875	1502	1511	1520	rVB4	72557	154268	1.06%	0.288%
48	12.034	1531	1538	1542	rBV	377283	619434	4.25%	1.154%
49	12.069	1542	1544	1549	rVB	143355	170596	1.17%	0.318%
50	12.128	1549	1554	1560	rBV2	67114	139851	0.96%	0.261%

51	12.275	1572	1579	1587	rVB2	288988	592133	4.06%	1.104%
52	12.439	1602	1607	1611	rBV3	119769	190878	1.31%	0.356%
53	12.487	1611	1615	1619	rVB	79578	117090	0.80%	0.218%
54	12.681	1641	1648	1661	rVB5	101032	279283	1.91%	0.521%
55	12.781	1661	1665	1672	rBV5	43152	78023	0.53%	0.145%

56	12.845	1672	1676	1682	rBV	58476	94958	0.65%	0.177%
57	12.992	1696	1701	1710	rVV2	110859	193535	1.33%	0.361%
58	13.281	1744	1750	1757	rVB	404965	594922	4.08%	1.109%
59	13.498	1780	1787	1794	rVB6	62401	153788	1.05%	0.287%
60	13.628	1800	1809	1814	rBV6	92900	232943	1.60%	0.434%

61	13.763	1827	1832	1836	rBV2	101870	199523	1.37%	0.372%
62	13.816	1836	1841	1844	rVV3	88338	172476	1.18%	0.321%
63	13.851	1844	1847	1853	rVB2	66183	106746	0.73%	0.199%
64	13.933	1858	1861	1865	rVB2	126834	176446	1.21%	0.329%
65	14.351	1927	1932	1937	rBV	2089846	2718649	18.63%	5.067%

66	14.445	1937	1948	1954	rVB	1561046	2388413	16.37%	4.451%
67	14.522	1958	1961	1967	rVB4	73827	107023	0.73%	0.199%
68	14.586	1968	1972	1979	rVB	321595	425337	2.92%	0.793%
69	14.657	1979	1984	1992	rVB7	80425	181739	1.25%	0.339%
70	14.845	2013	2016	2021	rVB4	81218	110507	0.76%	0.206%

71	14.910	2021	2027	2031	rBV4	64695	108883	0.75%	0.203%
72	14.975	2034	2038	2041	rBV4	75540	123882	0.85%	0.231%
73	15.063	2047	2053	2058	rBV	1163956	1641671	11.25%	3.060%
74	15.204	2073	2077	2085	rVB2	137329	242046	1.66%	0.451%
75	15.298	2085	2093	2098	rVB	462077	733461	5.03%	1.367%

76	15.433	2113	2116	2121	rVB	87931	110883	0.76%	0.207%
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Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

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 Title : SVOA CALIBRATION

77	15.539	2128	2134	2138	rBV	156000	235632	1.62%	0.439%
78	15.704	2157	2162	2167	rVB	140359	228386	1.57%	0.426%
79	15.798	2174	2178	2183	rVB4	93658	137830	0.94%	0.257%
80	15.916	2193	2198	2207	rVB6	63579	128948	0.88%	0.240%
81	16.163	2235	2240	2242	rBV	425542	583371	4.00%	1.087%
82	16.186	2242	2244	2248	rVV2	197636	217741	1.49%	0.406%
83	16.257	2252	2256	2260	rVV5	52353	79211	0.54%	0.148%
84	16.833	2344	2354	2367	rBV	10332348	14588942	100.00%	27.189%
85	16.951	2370	2374	2378	rVV	356549	488161	3.35%	0.910%
86	16.998	2378	2382	2391	rVB	171422	313894	2.15%	0.585%
87	17.180	2407	2413	2423	rVV	1959458	2934279	20.11%	5.469%
88	17.280	2426	2430	2434	rVV	106850	158393	1.09%	0.295%
89	17.322	2434	2437	2442	rVB2	102763	140286	0.96%	0.261%
90	17.474	2459	2463	2471	rVB5	79137	153741	1.05%	0.287%
91	17.686	2494	2499	2505	rVB	261658	323463	2.22%	0.603%
92	17.757	2507	2511	2516	rVV	123325	157710	1.08%	0.294%
93	18.374	2612	2616	2622	rBV	177014	226801	1.55%	0.423%
94	19.010	2720	2724	2731	rVB2	130737	236302	1.62%	0.440%
95	19.563	2814	2818	2820	rBV2	91400	132929	0.91%	0.248%
96	20.115	2908	2912	2918	rBV3	95454	170444	1.17%	0.318%
97	20.610	2993	2996	3000	rBV2	77681	93701	0.64%	0.175%
98	21.092	3074	3078	3083	rBV	183807	217568	1.49%	0.405%
99	21.404	3125	3131	3137	rBV	1899282	2893888	19.84%	5.393%
100	24.409	3634	3642	3656	rVB	1205784	3148793	21.58%	5.868%

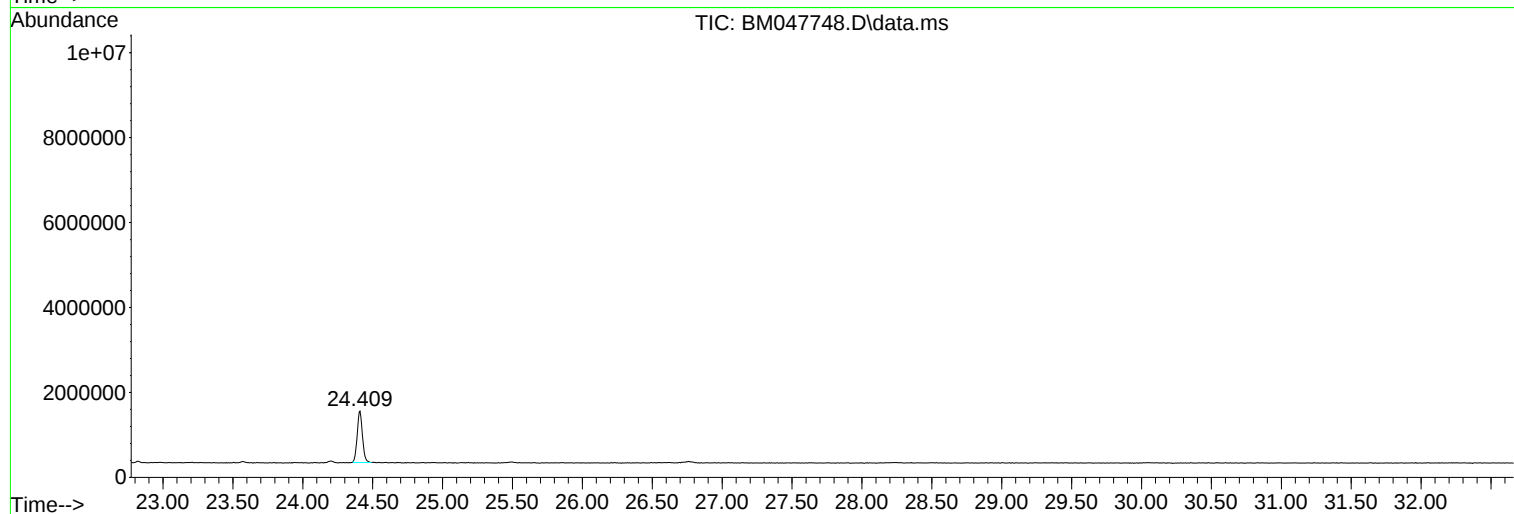
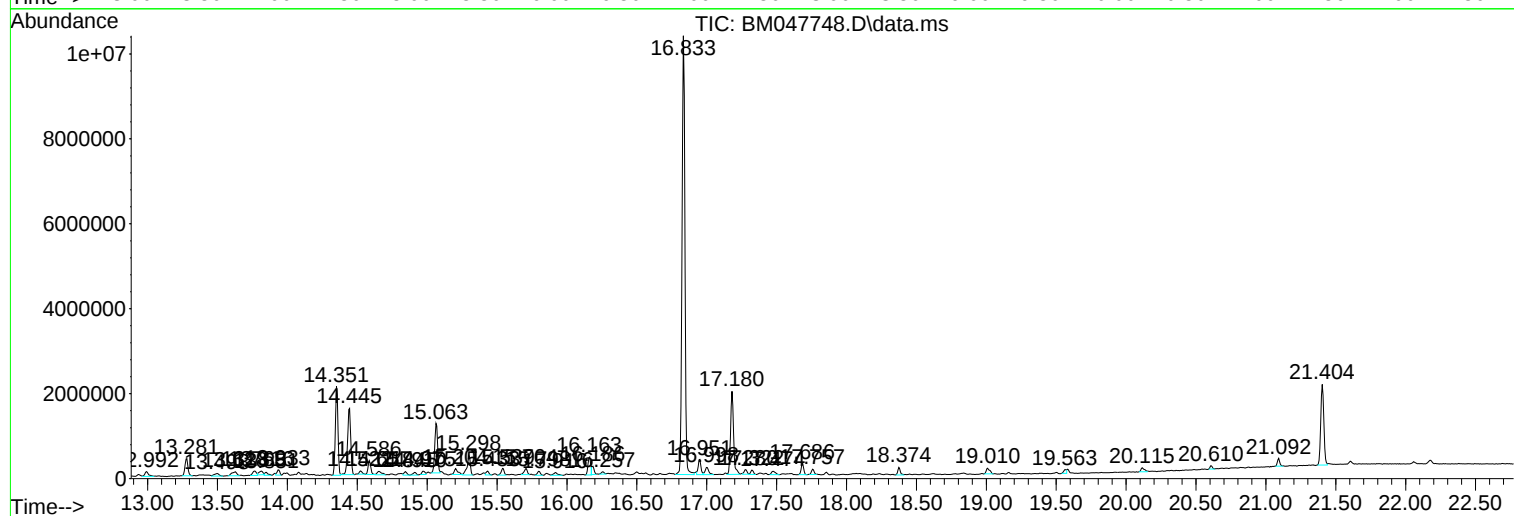
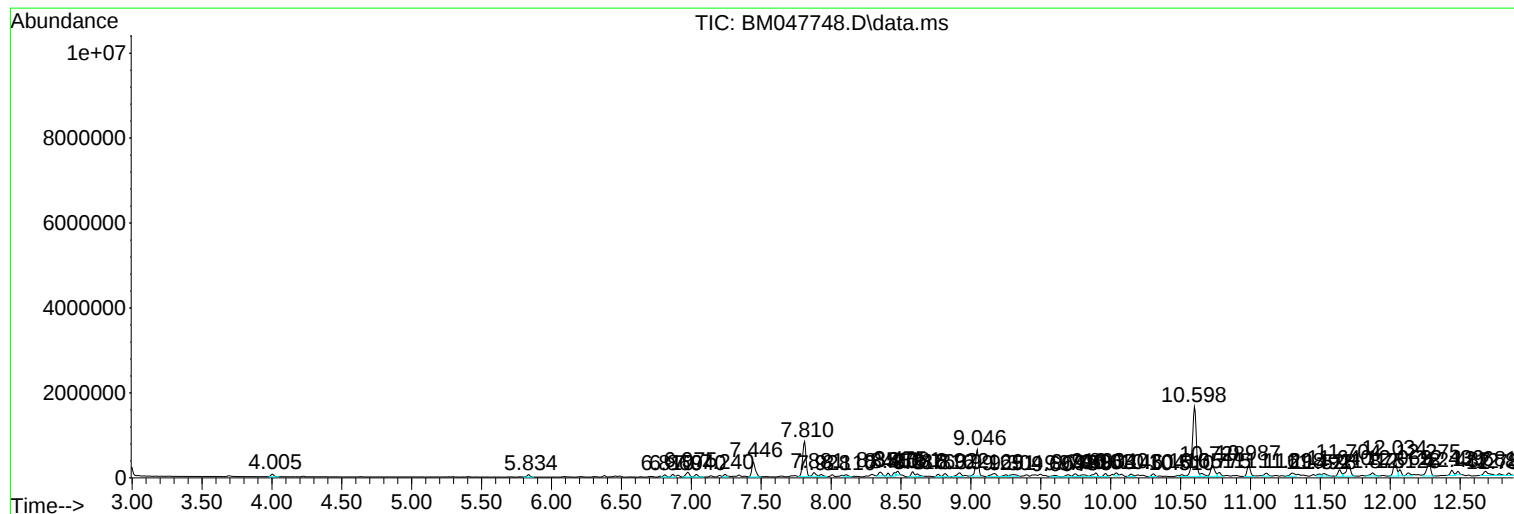
Sum of corrected areas: 53656597

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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCB4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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



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Acq On : 01 Oct 2024 18:21
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ALS Vial : 12 Sample Multiplier: 1

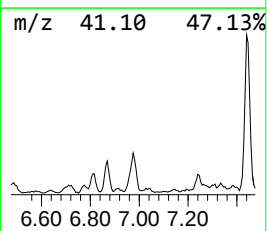
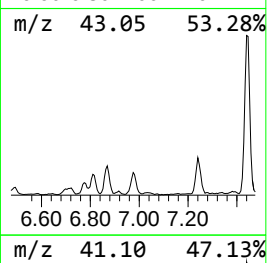
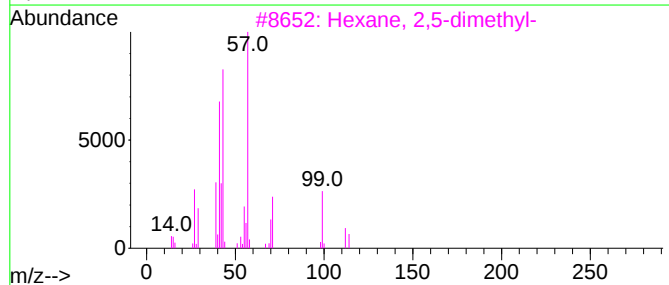
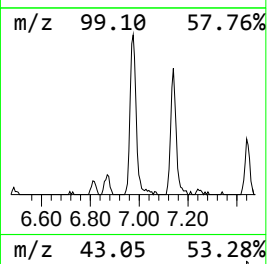
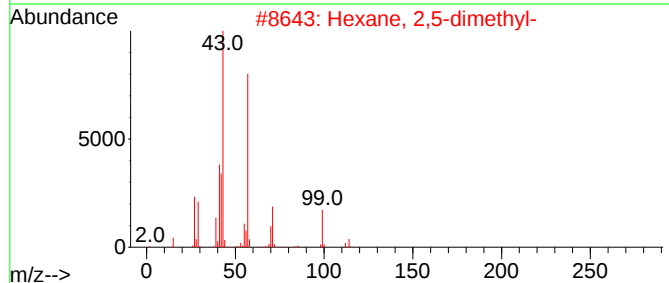
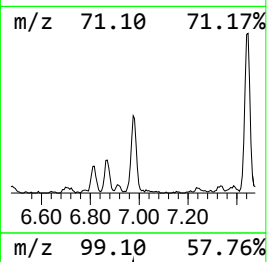
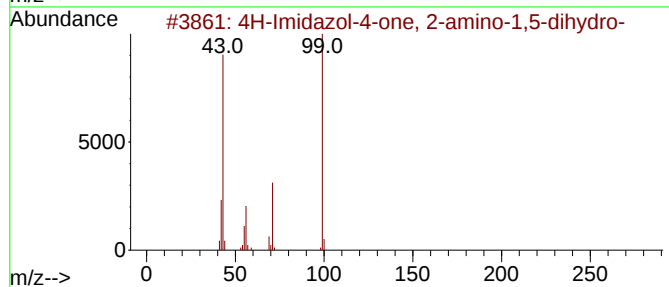
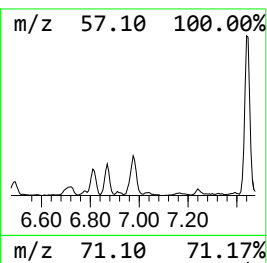
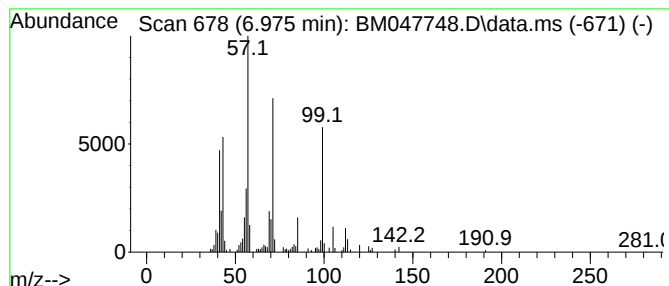
Instrument :
BNA_M
ClientSampleId :
DDCB4DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 4H-Imidazol-4-one, 2-amino-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
6.975	3.02 ng/ul	204385	1,4-Dichlorobenzene-d4	7.810	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4H-Imidazol-4-one, 2-amino-1,5-d...	99	C3H5N3O	000503-86-6	58
2	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	50
3	Hexane, 2,5-dimethyl-	114	C8H18	000592-13-2	47
4	3-Octanone, 2-methyl-	142	C9H18O	000923-28-4	43
5	Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	43



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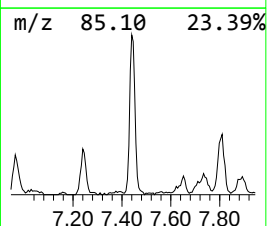
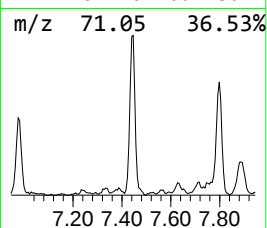
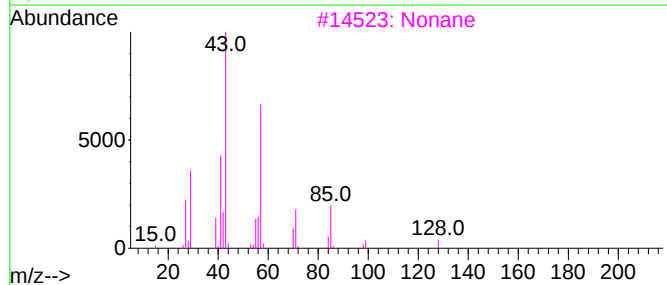
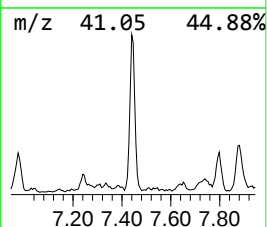
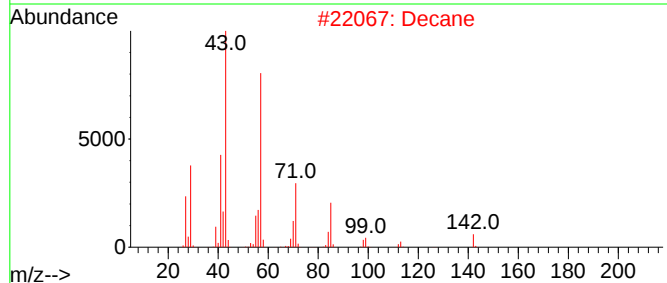
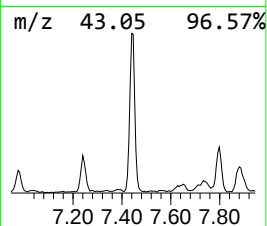
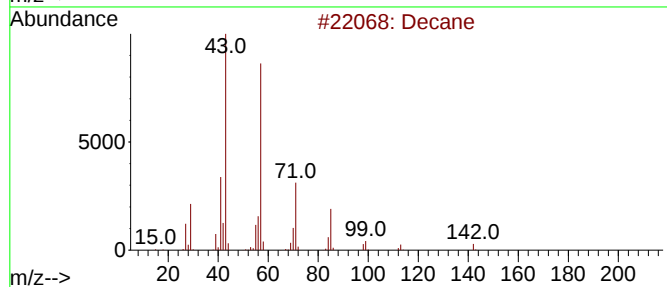
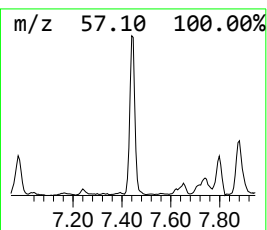
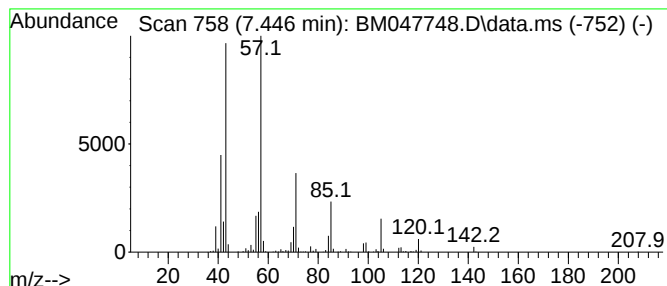
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.446	9.39 ng/ul	635597	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Decane	142	C10H22	000124-18-5	94
2			Decane	142	C10H22	000124-18-5	80
3			Nonane	128	C9H20	000111-84-2	80
4			Octane, 4-ethyl-	142	C10H22	015869-86-0	68
5			Carbonic acid, decyl vinyl ester	228	C13H24O3	1000383-25-7	64



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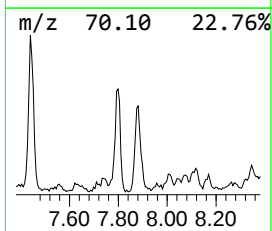
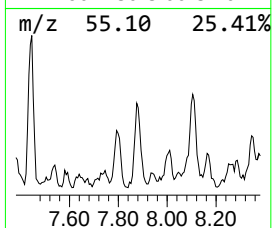
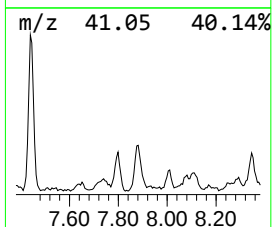
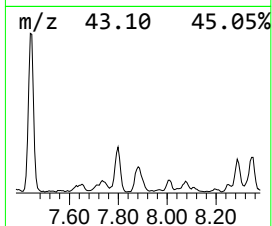
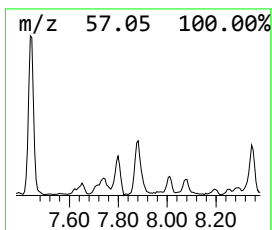
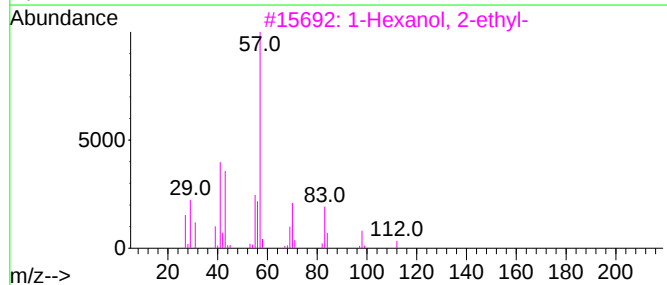
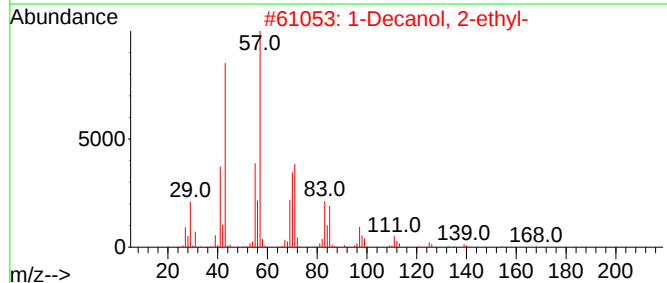
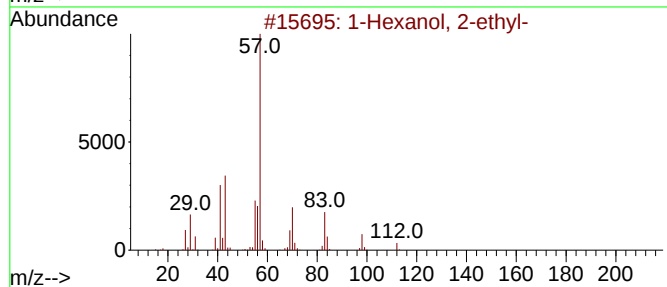
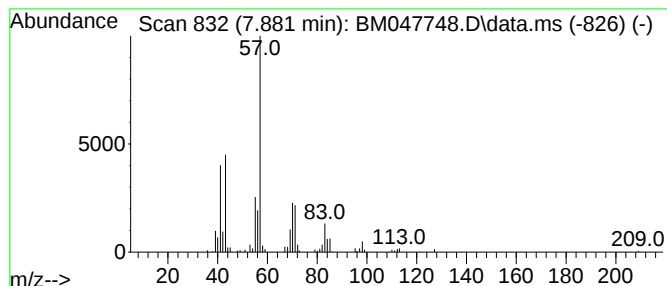
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 1-Hexanol, 2-ethyl- Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.881	2.16 ng/ul	145885	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	59
2	1-Decanol, 2-ethyl-			186	C12H26O	021078-65-9	53
3	1-Hexanol, 2-ethyl-			130	C8H18O	000104-76-7	53
4	2-Ethylhexyl mercaptoacetate			204	C10H20O2S	007659-86-1	50
5	Octane, 2,3-dimethyl-			142	C10H22	007146-60-3	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

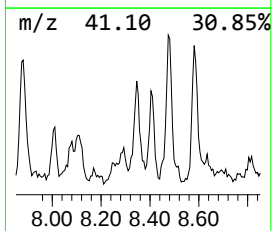
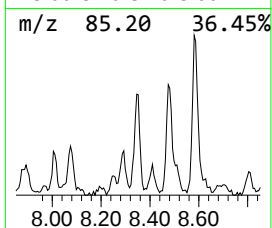
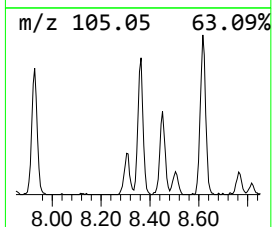
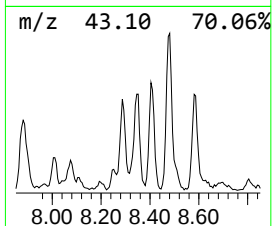
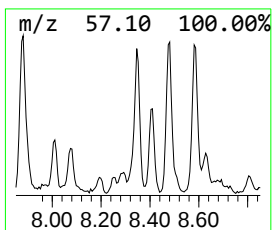
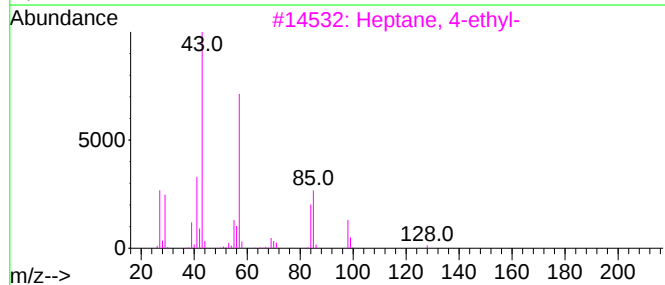
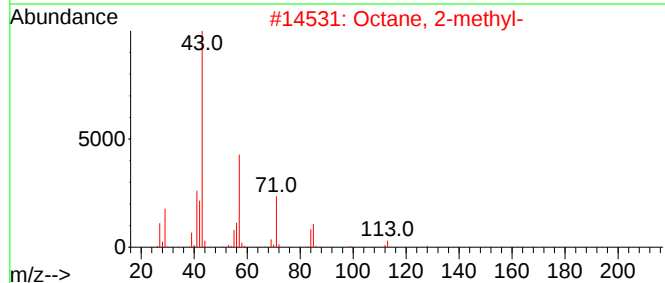
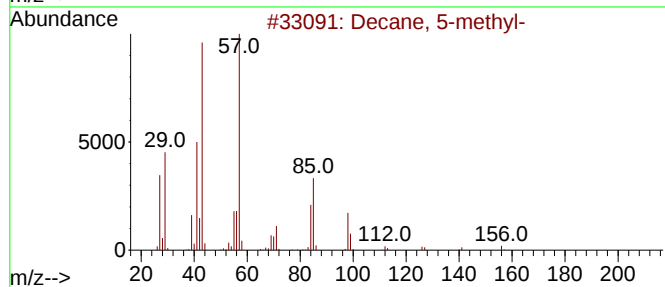
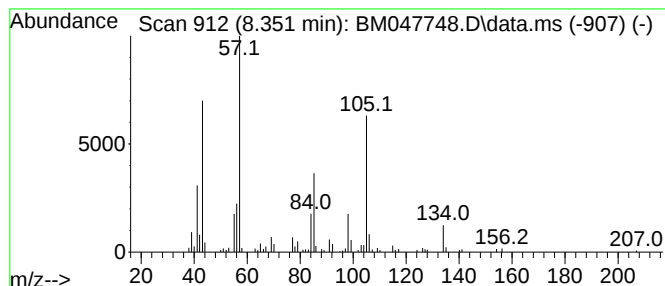
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD			R.T.
8.351	2.42 ng/ul	163555	1,4-Dichlorobenzene-d4			7.810
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Decane, 5-methyl-		156	C11H24	013151-35-4	58
2	Octane, 2-methyl-		128	C9H20	003221-61-2	38
3	Heptane, 4-ethyl-		128	C9H20	002216-32-2	38
4	Hexane, 2,3,5-trimethyl-		128	C9H20	001069-53-0	38
5	Heptane, 2,3-dimethyl-		128	C9H20	003074-71-3	27



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
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Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

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ClientSampleId :
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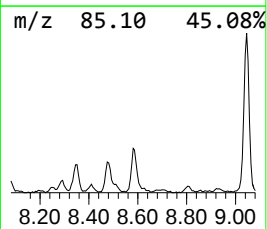
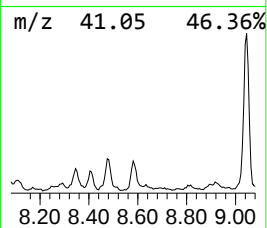
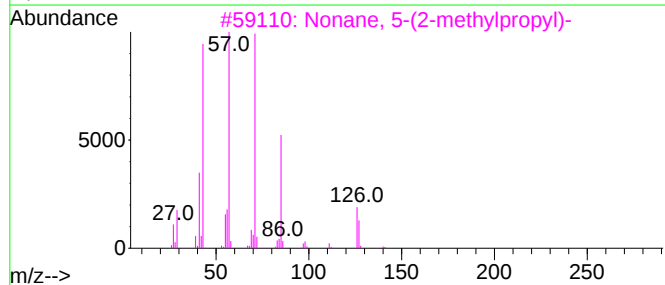
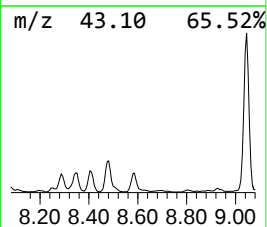
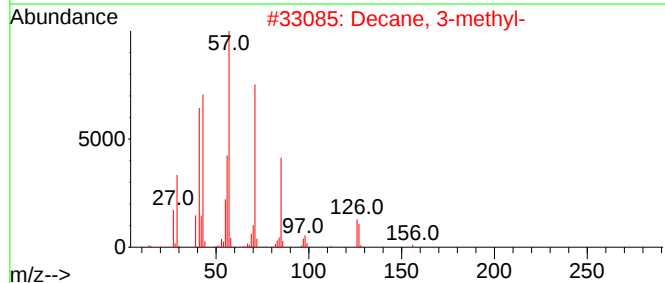
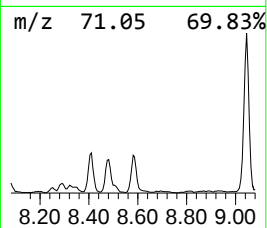
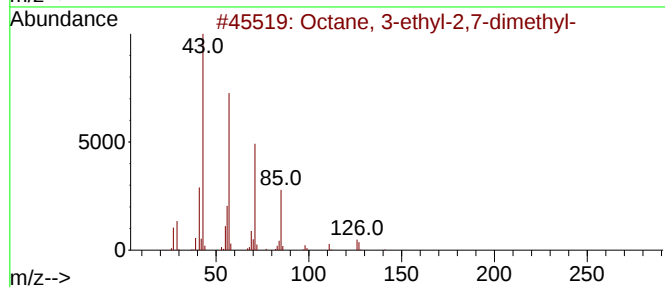
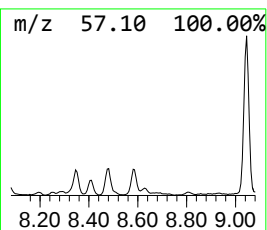
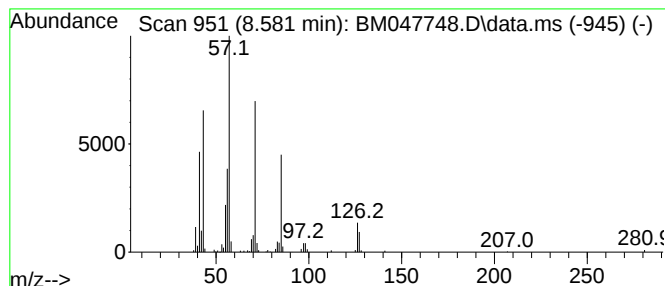
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.581	2.87 ng/ul	194165	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octane, 3-ethyl-2,7-dimethyl-	170	C12H26	062183-55-5	83		
2	Decane, 3-methyl-	156	C11H24	013151-34-3	80		
3	Nonane, 5-(2-methylpropyl)-	184	C13H28	062185-53-9	72		
4	Nonane, 5-butyl-	184	C13H28	017312-63-9	59		
5	2-Bromo dodecane	248	C12H25Br	013187-99-0	59		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
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Sample : P4018-08DL 30X
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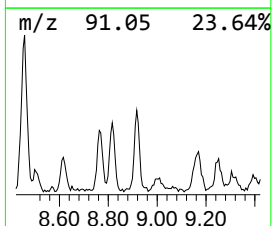
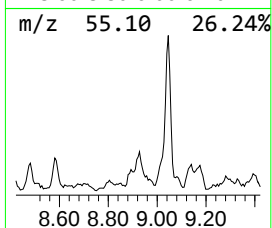
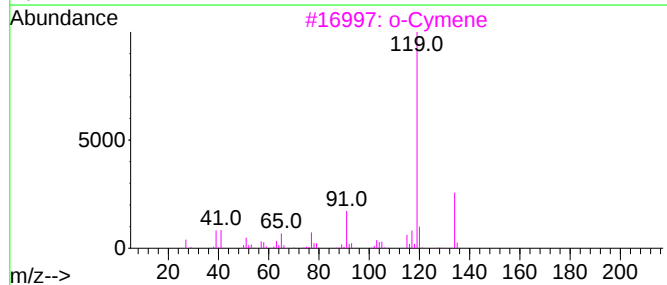
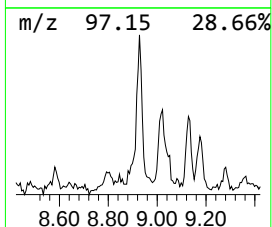
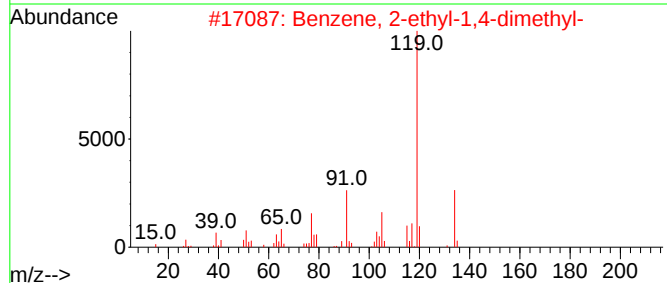
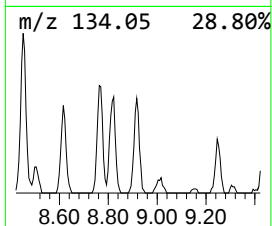
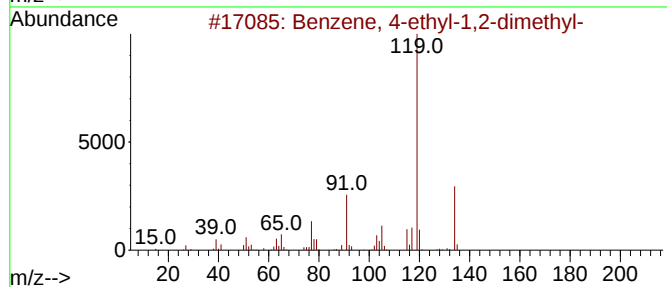
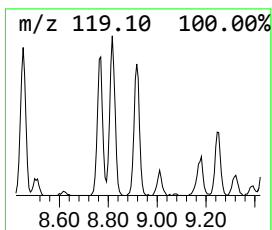
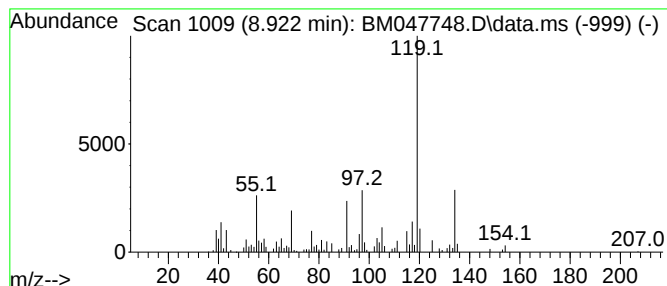
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.922	2.95 ng/ul	199492	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
3			o-Cymene	134	C10H14	000527-84-4	94
4			p-Cymene	134	C10H14	000099-87-6	93
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
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Instrument :
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ClientSampleId :
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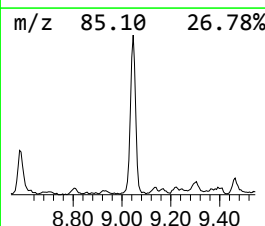
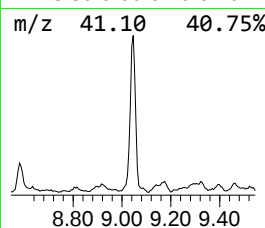
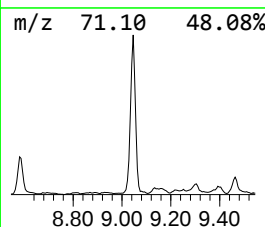
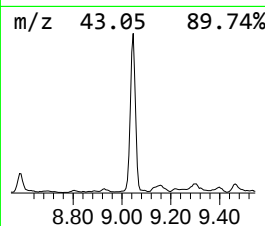
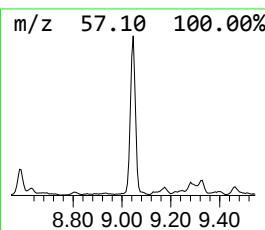
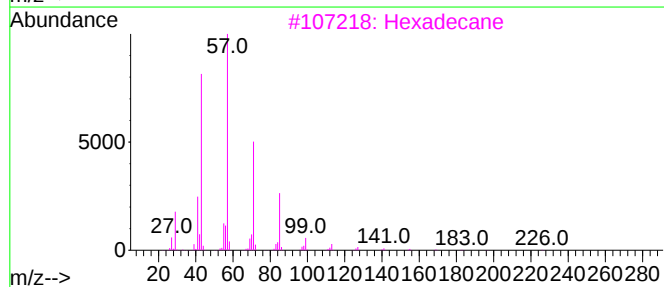
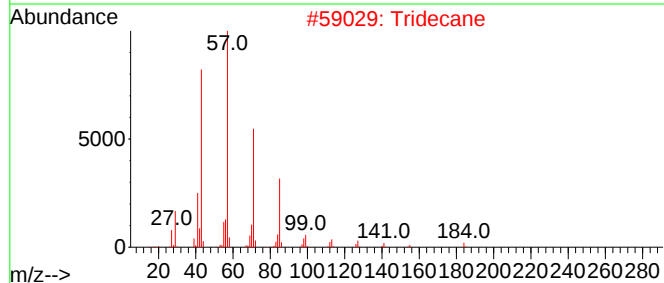
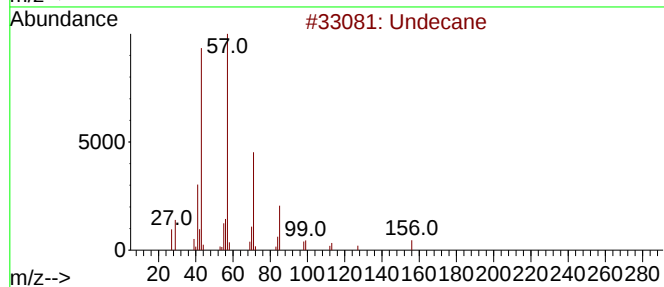
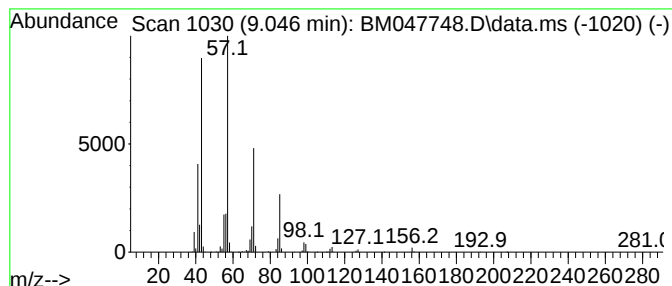
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.046	13.58 ng/ul	919055	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Undecane	156	C11H24	001120-21-4	90
2			Tridecane	184	C13H28	000629-50-5	90
3			Hexadecane	226	C16H34	000544-76-3	86
4			Decane, 2-methyl-	156	C11H24	006975-98-0	80
5			Tridecane	184	C13H28	000629-50-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
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Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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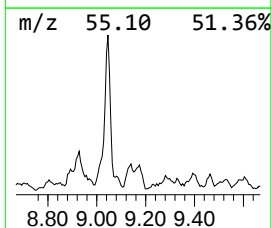
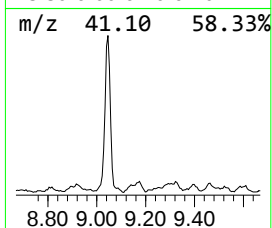
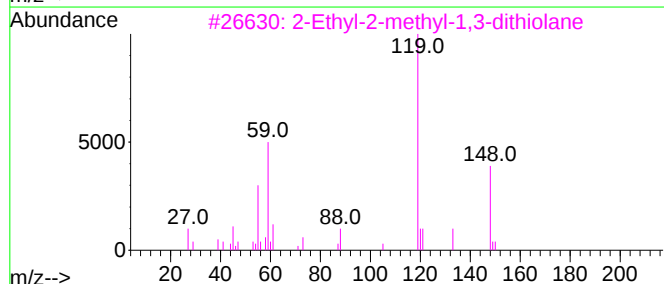
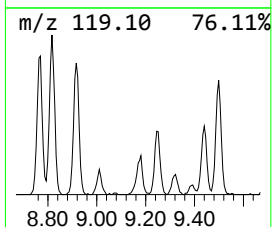
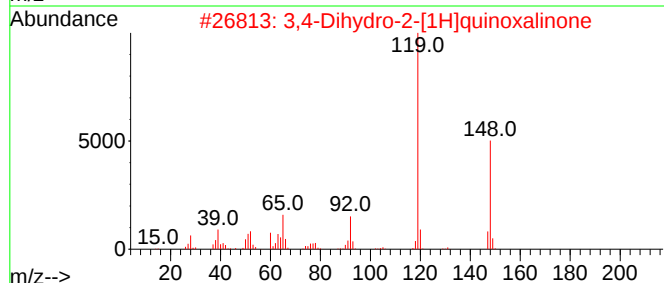
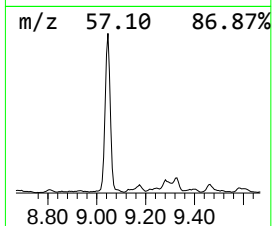
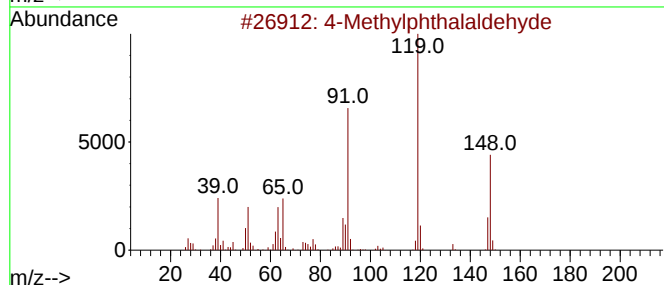
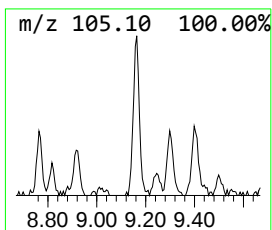
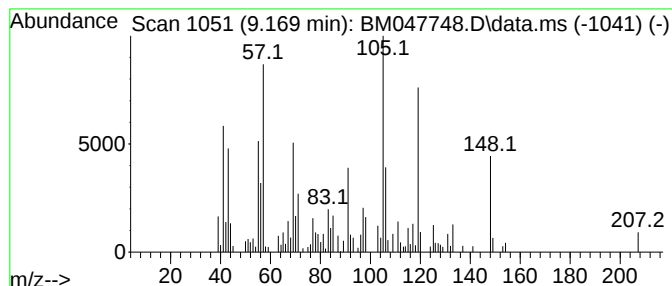
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-01 Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.169	3.01 ng/ul	203812	1,4-Dichlorobenzene-d4	7.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-Methylphthalaldehyde	148	C9H8O2	015158-36-8	49
2			3,4-Dihydro-2-[1H]quinoxalinone	148	C8H8N2O	059564-59-9	43
3			2-Ethyl-2-methyl-1,3-dithiolane	148	C6H12S2	006008-81-7	43
4			Pyridine, 5-ethenyl-2-methyl-	119	C8H9N	000140-76-1	30
5			4-Aminostyrene	119	C8H9N	001520-21-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
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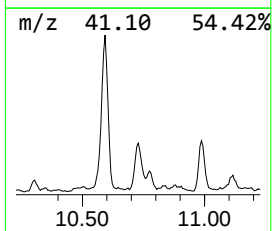
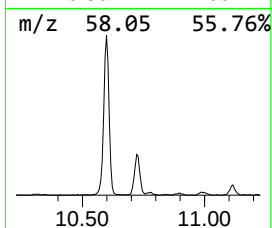
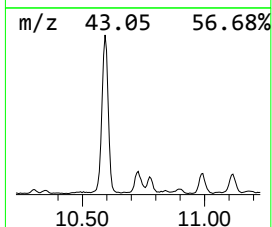
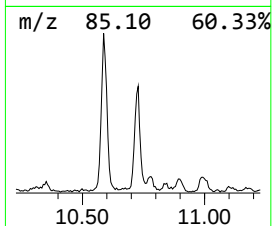
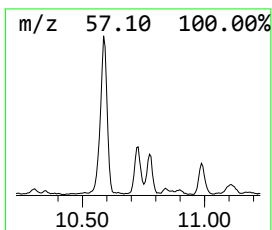
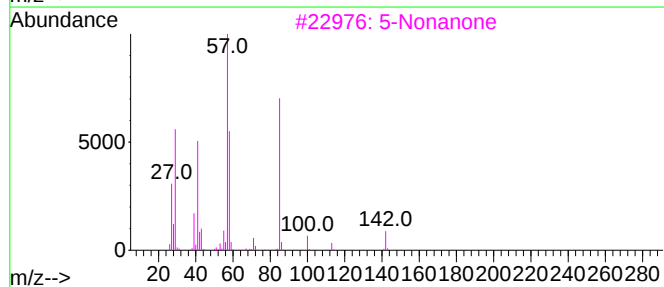
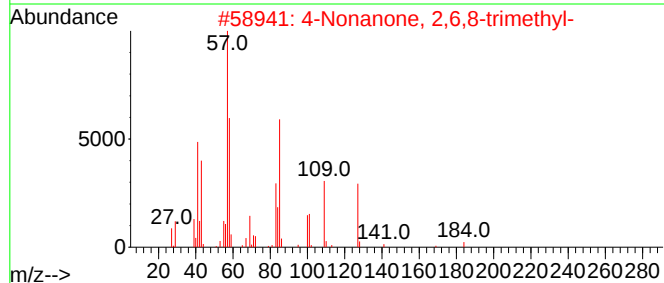
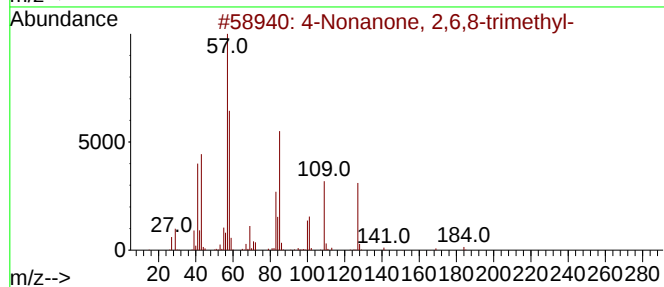
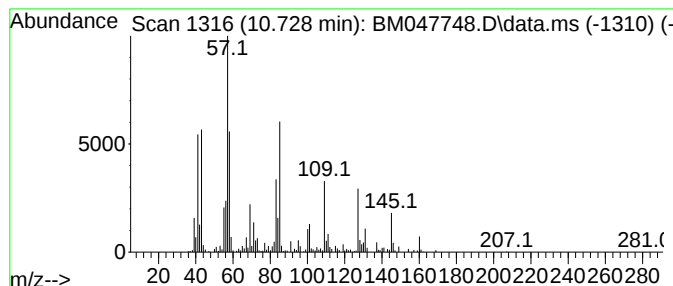
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 4-Nonanone, 2,6,8-trimethyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
10.728	3.26 ng/ul	521406	Naphthalene-d8	10.604	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	86
2	4-Nonanone, 2,6,8-trimethyl-	184	C12H24O	000123-18-2	86
3	5-Nonanone	142	C9H18O	000502-56-7	35
4	5-Nonanone	142	C9H18O	000502-56-7	35
5	Thiazole	85	C3H3NS	000288-47-1	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
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Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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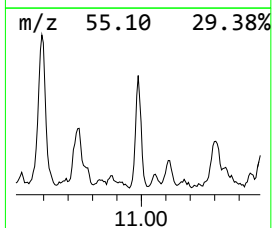
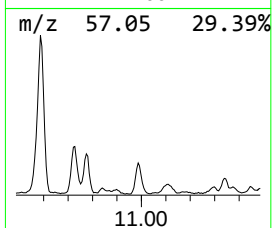
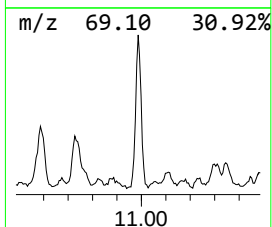
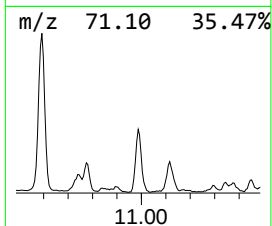
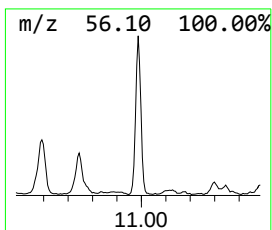
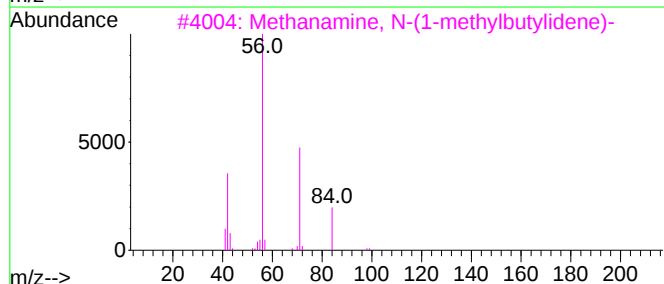
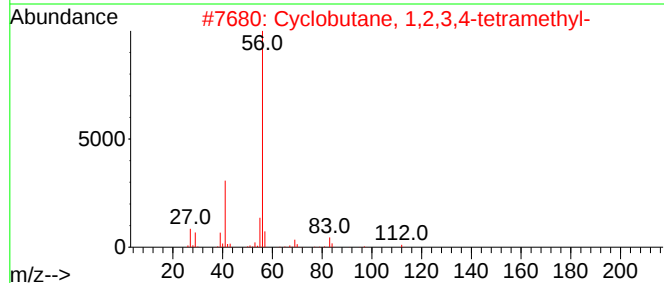
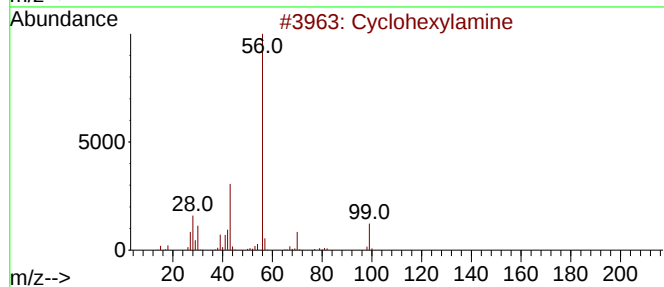
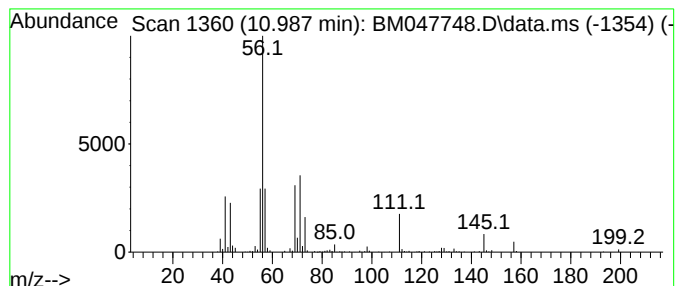
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-02 Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.987	2.83 ng/ul	454061	Naphthalene-d8	10.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexylamine	99	C6H13N	000108-91-8	43
2			Cyclobutane, 1,2,3,4-tetramethyl-	112	C8H16	069531-57-3	43
3			Methanamine, N-(1-methylbutylidene)-	99	C6H13N	022431-09-0	42
4			2,4-Dipropyl-5-ethyl-1,3-dioxane	200	C12H24O2	006413-83-8	40
5			Tridecane, 7-methylene-	196	C14H28	019780-80-4	33



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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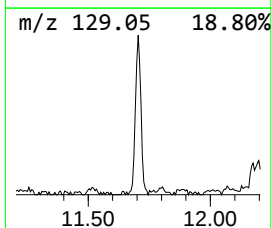
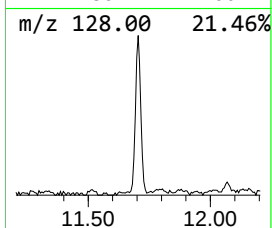
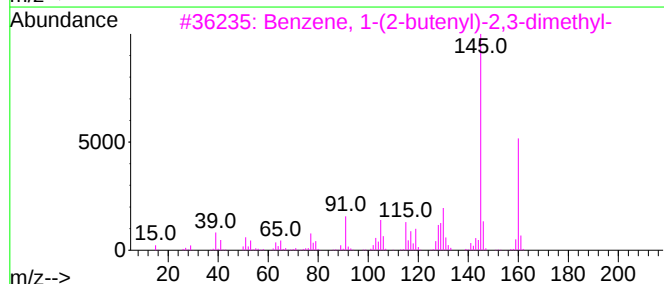
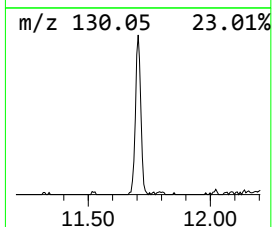
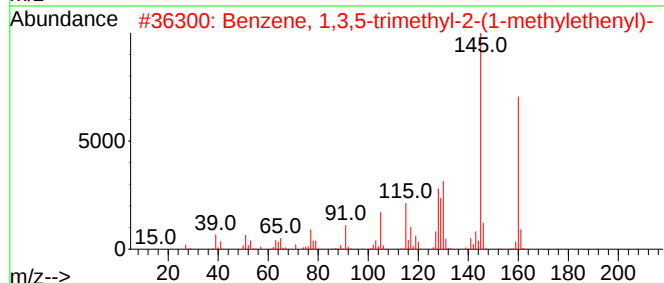
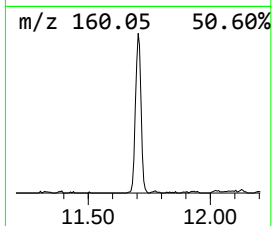
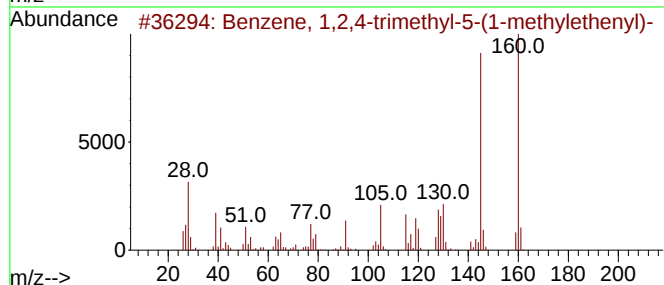
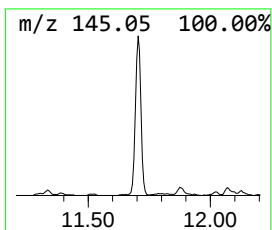
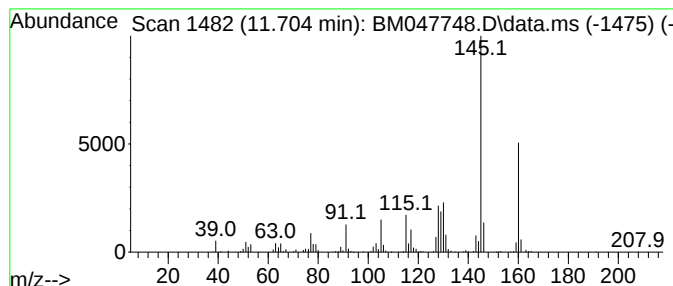
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 1,2,4-trimethyl-5-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.704	3.54 ng/ul	566992	Naphthalene-d8	10.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	95
2			Benzene, 1,3,5-trimethyl-2-(1-me...	160	C12H16	014679-13-1	94
3			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
4			Benzene, 1-(2-butenyl)-2,3-dimet...	160	C12H16	054340-85-1	94
5			Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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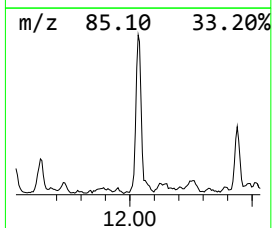
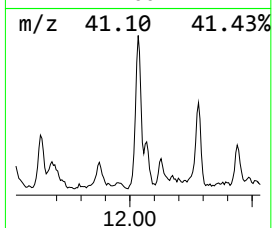
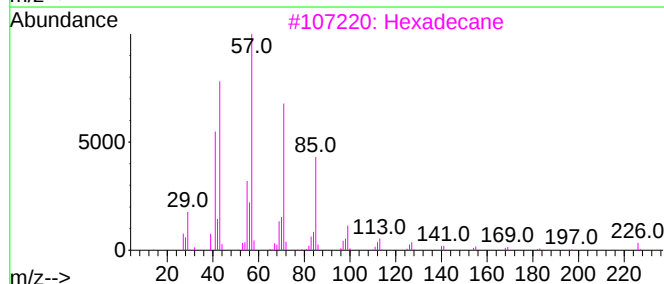
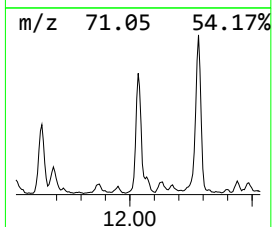
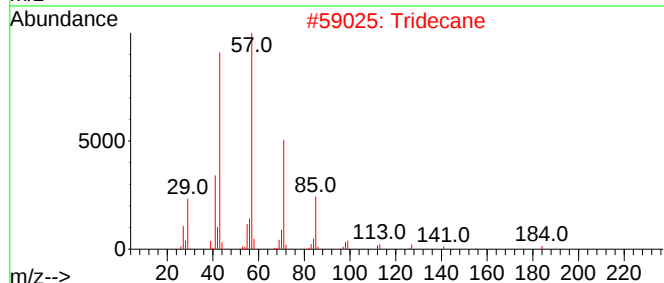
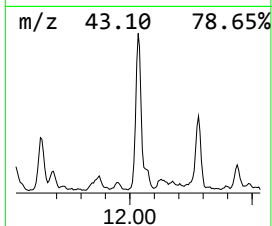
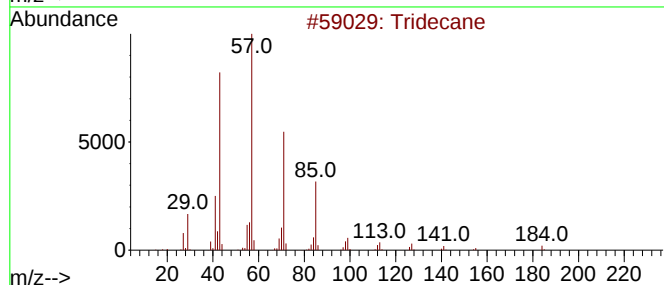
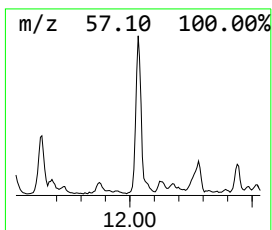
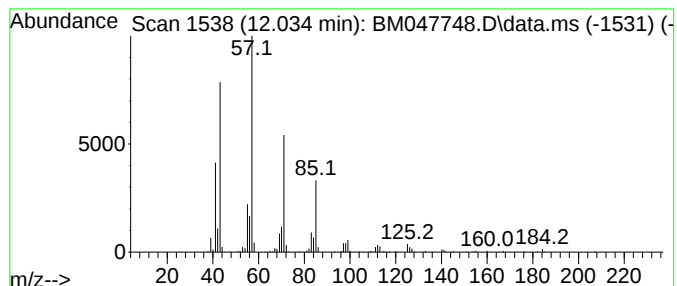
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.034	3.87 ng/ul	619434	Naphthalene-d8	10.604

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane	184	C13H28	000629-50-5	87
2			Tridecane	184	C13H28	000629-50-5	87
3			Hexadecane	226	C16H34	000544-76-3	80
4			Dodecane, 2-methyl-	184	C13H28	001560-97-0	74
5			Hexadecane	226	C16H34	000544-76-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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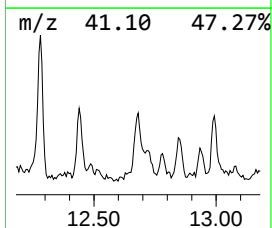
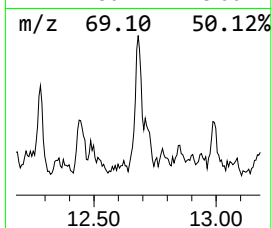
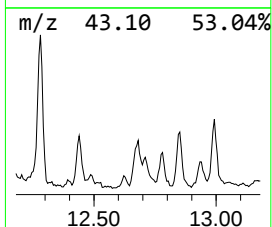
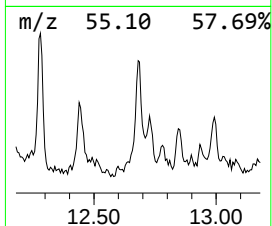
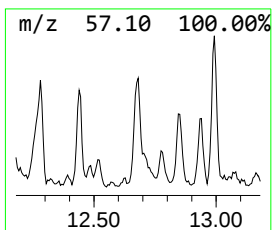
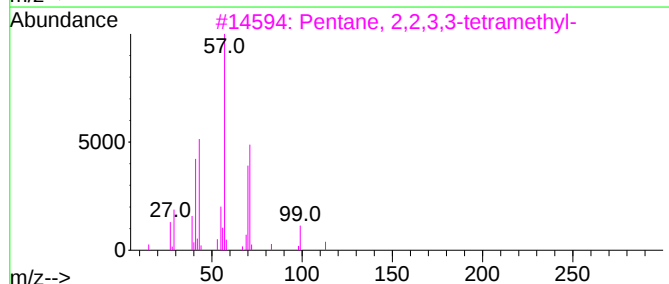
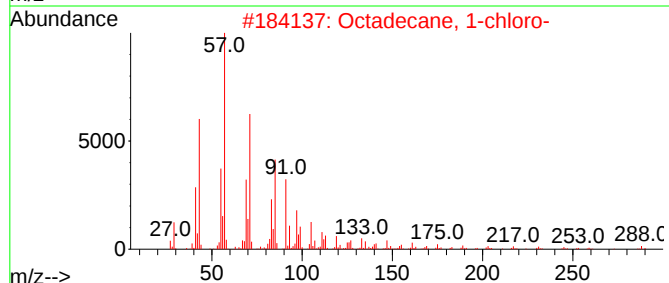
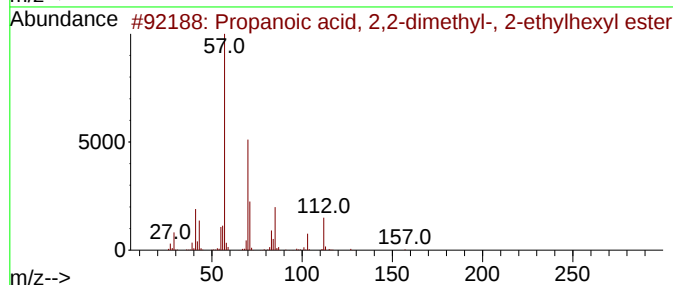
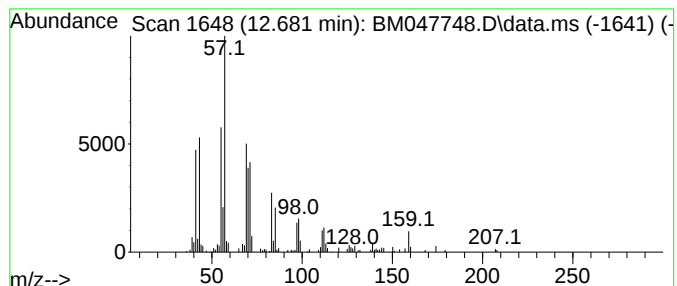
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 unknown-03 Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.681	2.34 ng/ul	279283	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Propanoic acid, 2,2-dimethyl-, 2...	214	C13H26O2	016387-18-1	43
2			Octadecane, 1-chloro-	288	C18H37Cl	003386-33-2	43
3			Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	38
4			Undecane, 3,9-dimethyl-	184	C13H28	017301-31-4	35
5			Nonane, 3,7-dimethyl-	156	C11H24	017302-32-8	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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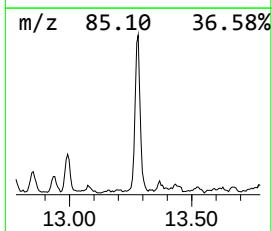
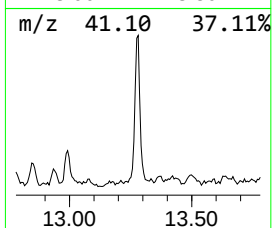
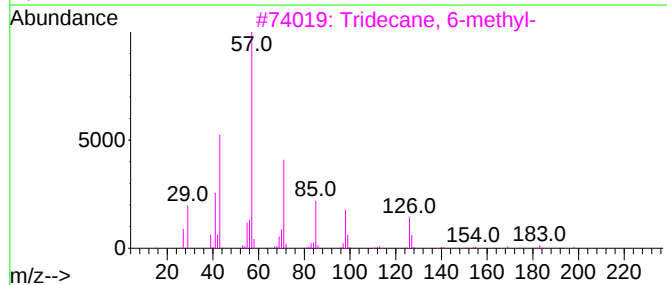
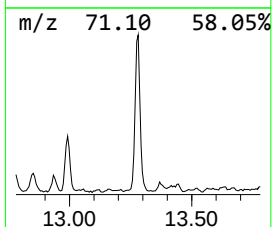
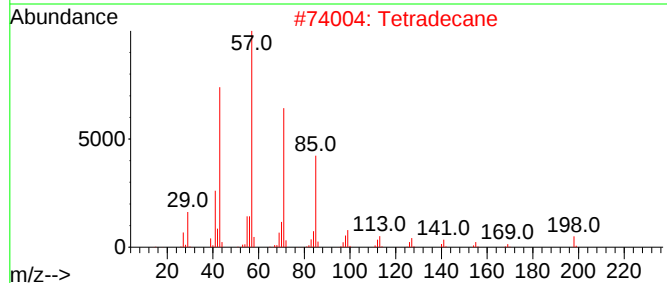
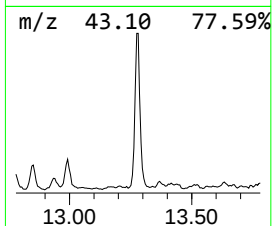
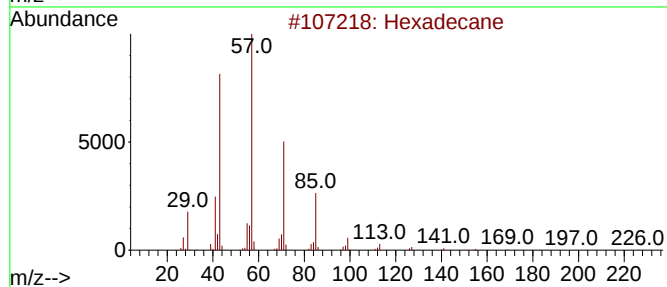
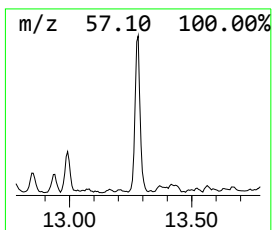
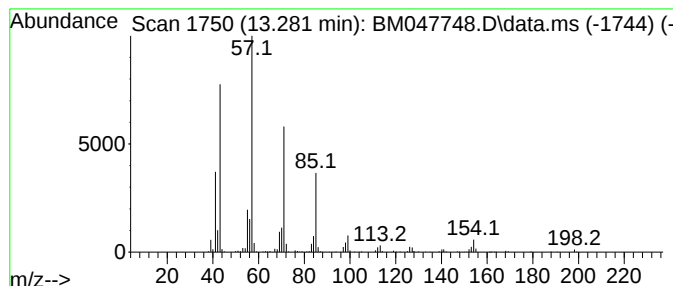
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.281	4.98 ng/ul	594922	Acenaphthene-d10	14.445

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane	226	C16H34	000544-76-3	87
2		Tetradecane	198	C14H30	000629-59-4	87
3		Tridecane, 6-methyl-	198	C14H30	013287-21-3	87
4		Hexadecane	226	C16H34	000544-76-3	86
5		Tridecane	184	C13H28	000629-50-5	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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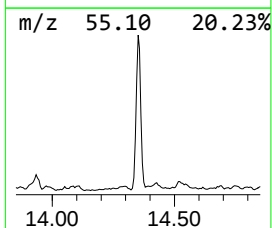
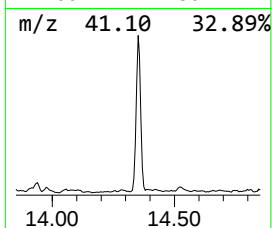
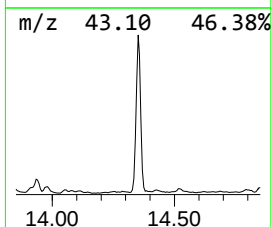
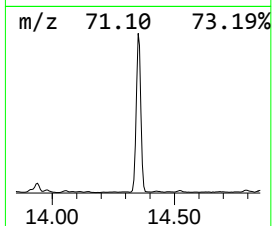
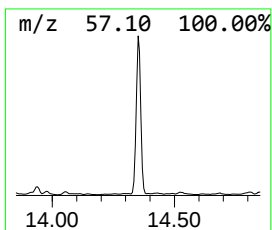
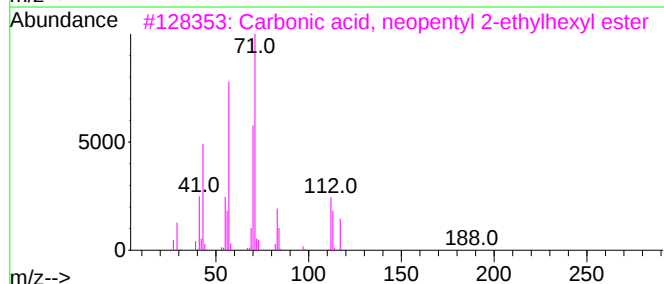
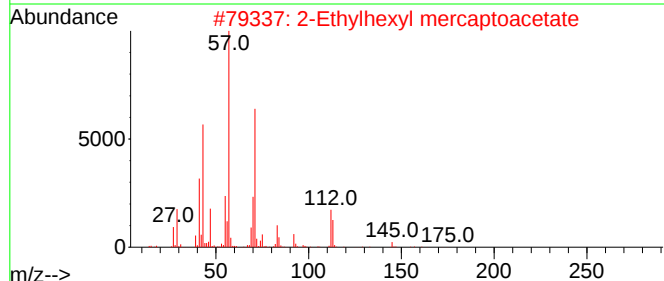
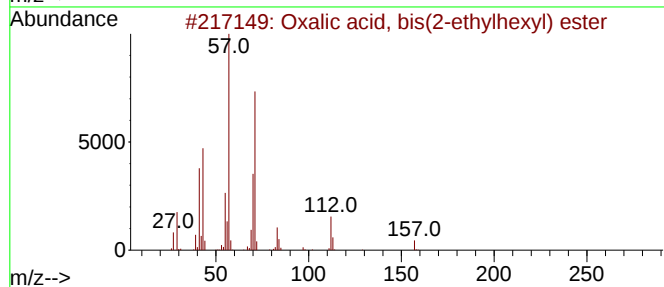
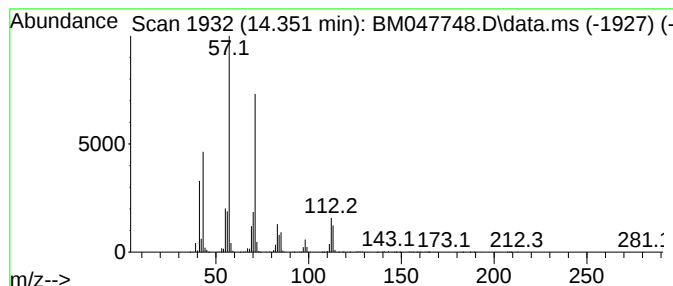
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Oxalic acid, bis(2-ethylhex... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.351	22.77 ng/ul	2718650	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxalic acid, bis(2-ethylhexyl) e...	314	C18H34O4	1000309-39-0	80
2			2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	72
3			Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	72
4			Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	72
5			Sulfurous acid, decyl 2-ethylhex...	334	C18H38O3S	1000309-19-3	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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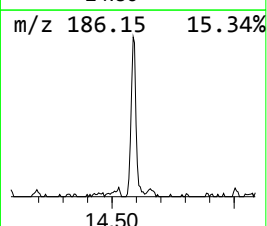
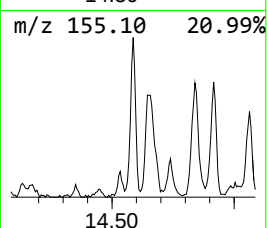
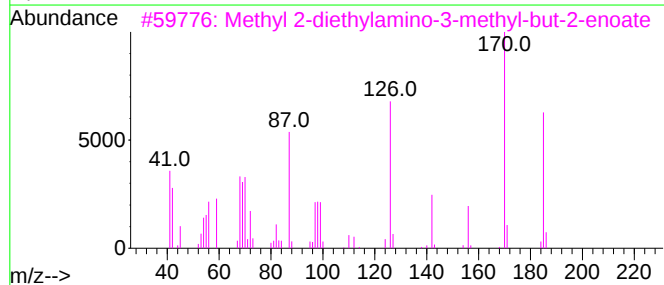
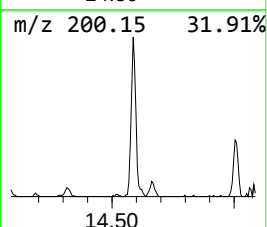
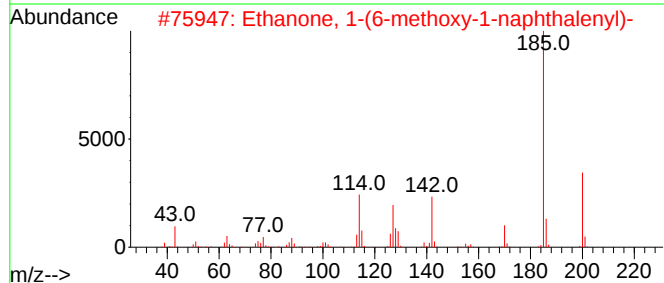
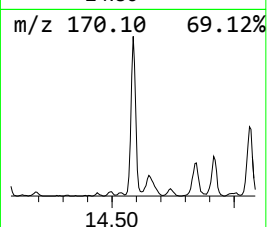
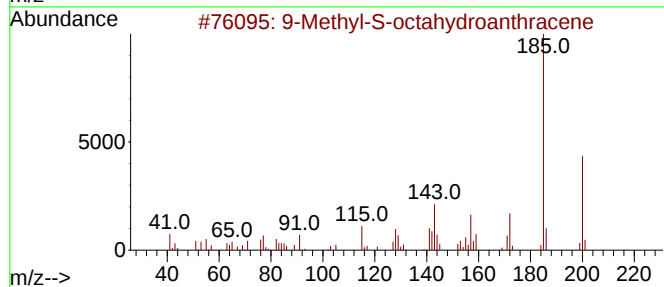
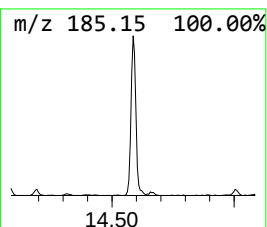
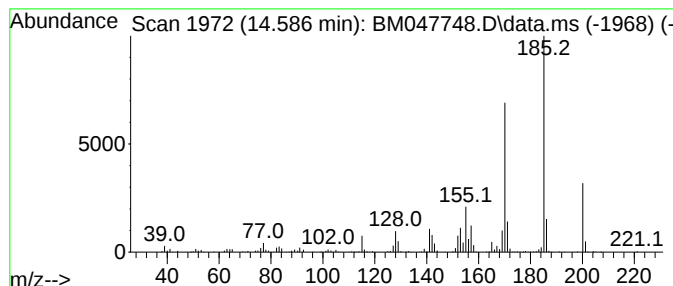
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 9-Methyl-S-octahydroanthracene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.586	3.56 ng/ul	425337	Acenaphthene-d10	14.445

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			9-Methyl-S-octahydroanthracene	200	C15H20	004948-51-0	60
2			Ethanone, 1-(6-methoxy-1-naphtha...	200	C13H12O2	058149-89-6	47
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_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

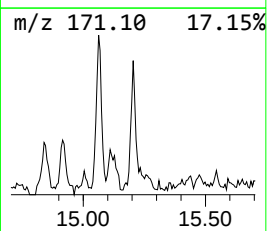
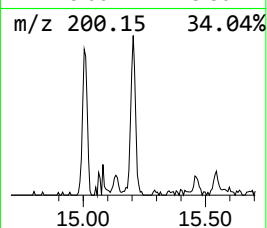
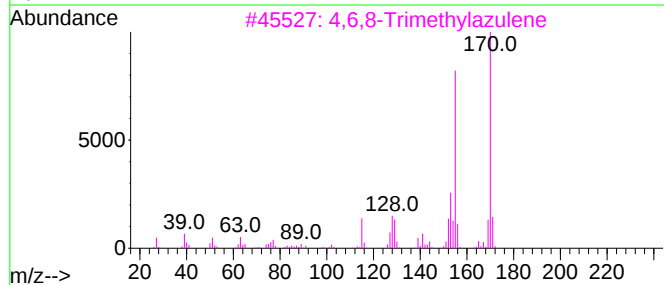
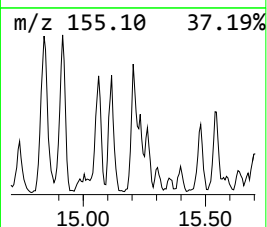
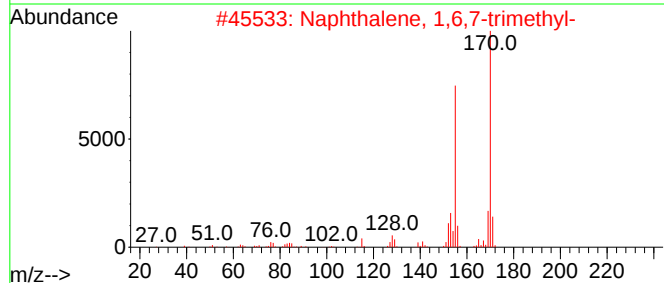
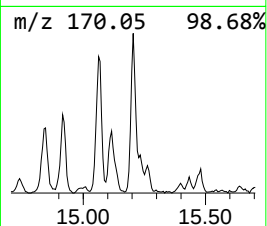
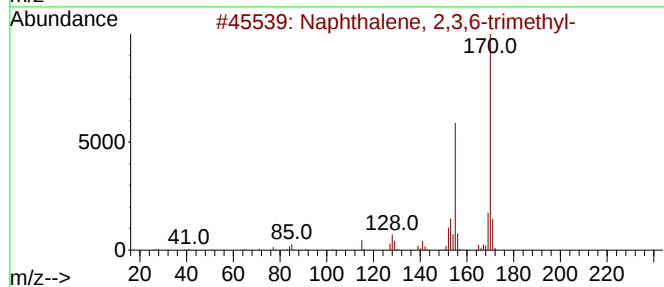
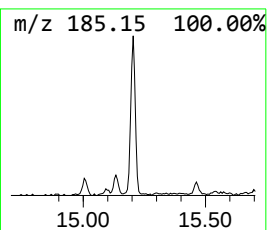
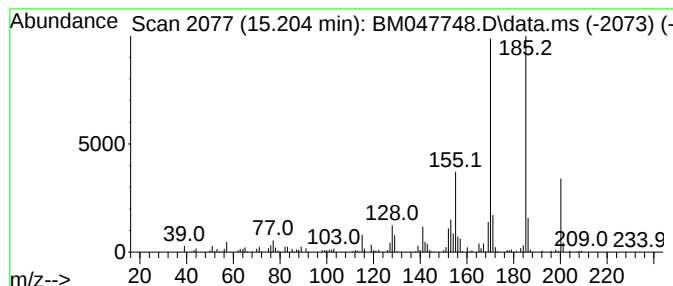
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 Naphthalene, 2,3,6-trimethyl- Concentration Rank 22

R.T.	EstConc	Area	Relative to ISTD	R.T.	
15.204	2.03 ng/ul	242046	Acenaphthene-d10	14.445	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2,3,6-trimethyl-	170	C13H14	000829-26-5	87
2	Naphthalene, 1,6,7-trimethyl-	170	C13H14	002245-38-7	70
3	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	70
4	4,6,8-Trimethylazulene	170	C13H14	000941-81-1	55
5	Methyl 2-diethylamino-3-methyl-b...	185	C10H19NO2	075122-81-5	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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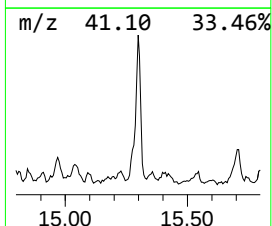
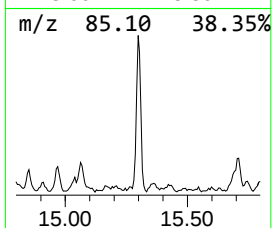
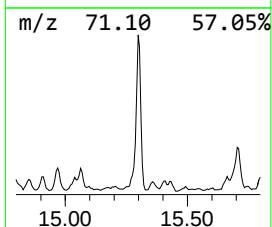
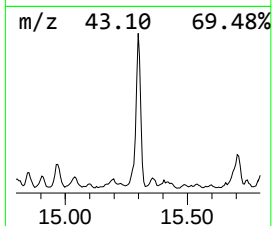
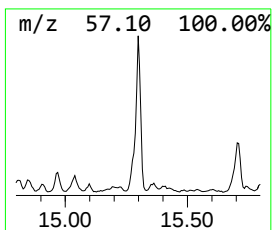
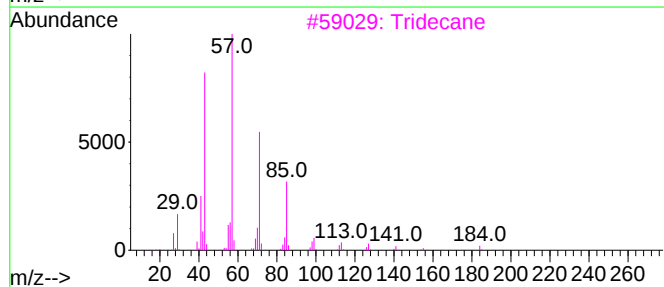
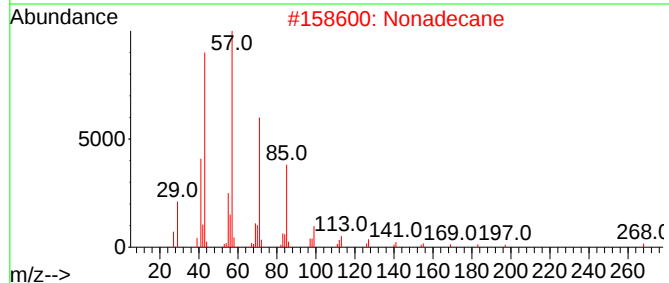
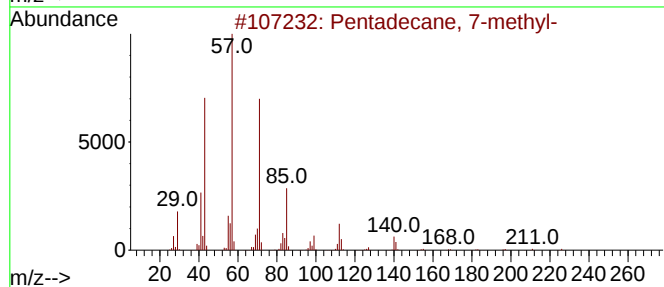
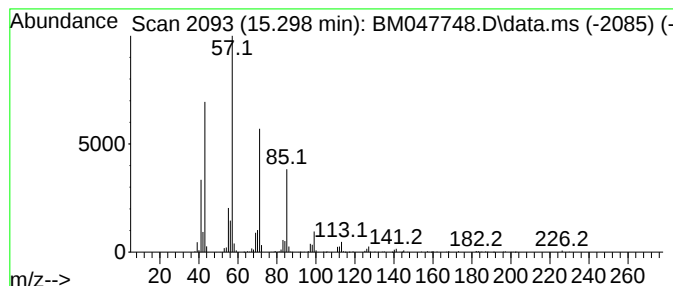
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.298	6.14 ng/ul	733461	Acenaphthene-d10	14.445

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Pentadecane, 7-methyl-	226	C16H34	006165-40-8	91
2		Nonadecane	268	C19H40	000629-92-5	91
3		Tridecane	184	C13H28	000629-50-5	90
4		Pentadecane	212	C15H32	000629-62-9	90
5		Heptadecane	240	C17H36	000629-78-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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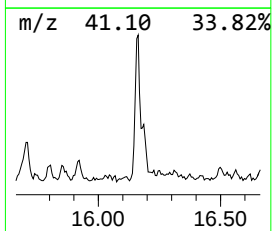
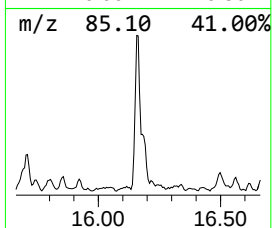
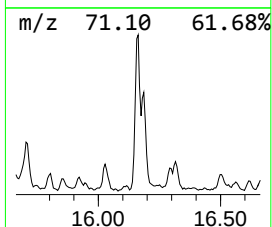
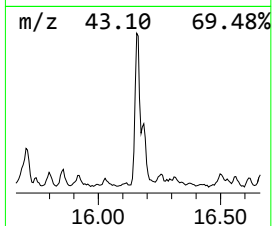
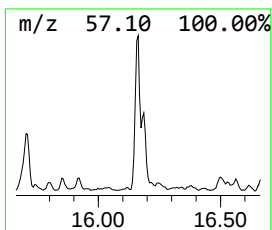
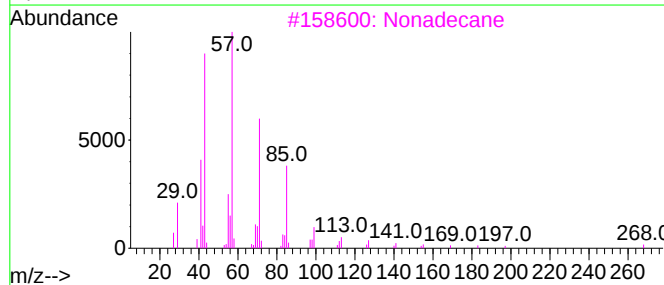
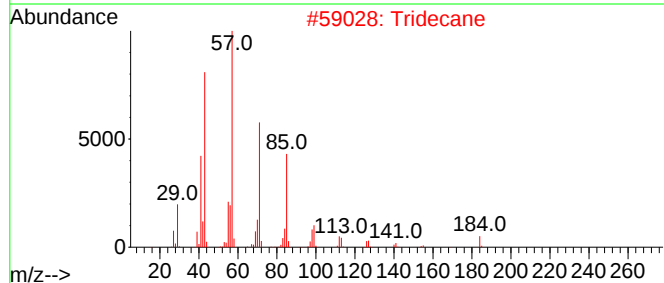
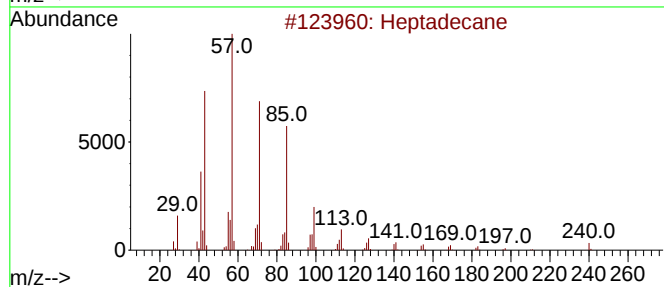
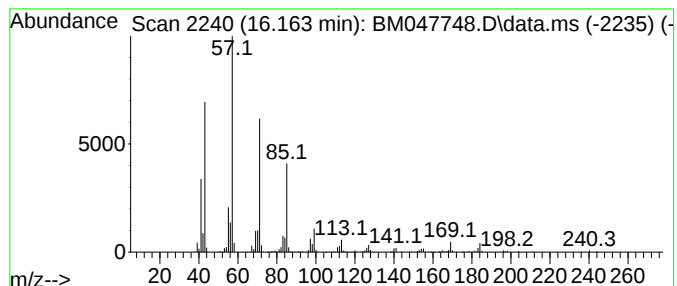
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.163	3.98 ng/ul	583371	Phenanthrene-d10	17.180

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	95
2		Tridecane	184	C13H28	000629-50-5	93
3		Nonadecane	268	C19H40	000629-92-5	91
4		Eicosane	282	C20H42	000112-95-8	87
5		Heptacosane	380	C27H56	000593-49-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
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Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

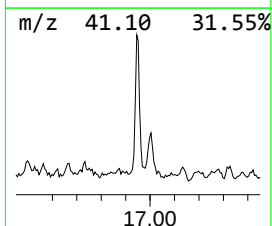
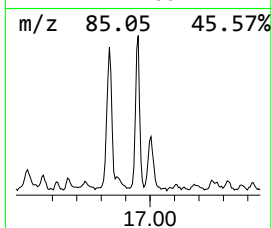
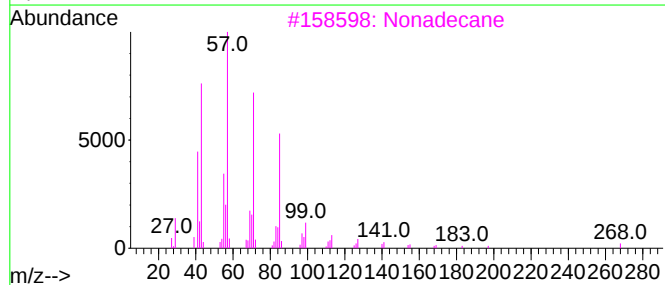
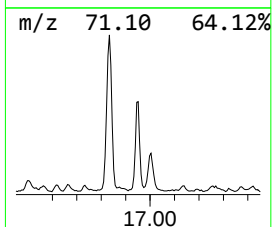
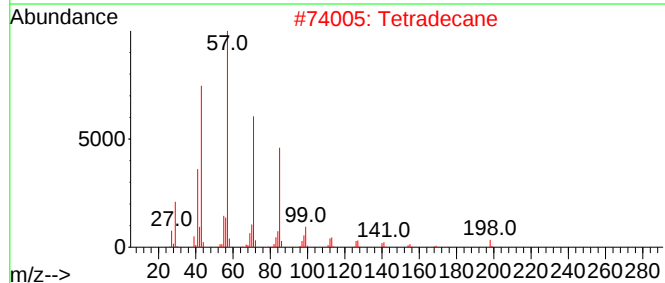
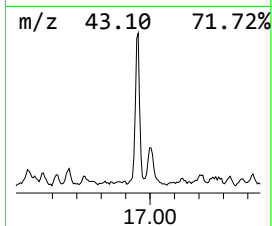
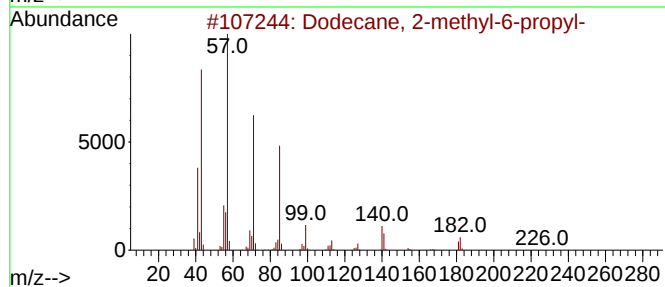
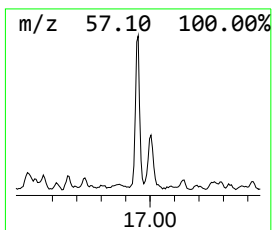
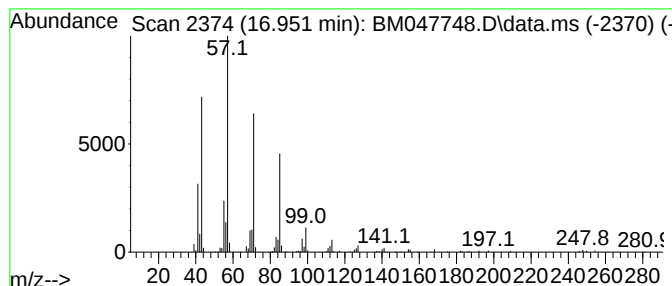
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD		R.T.	
16.951	3.33 ng/ul	488161	Phenanthrene-d10		17.180	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	86
2	Tetradecane		198	C14H30	000629-59-4	86
3	Nonadecane		268	C19H40	000629-92-5	86
4	Heneicosane		296	C21H44	000629-94-7	72
5	Decane, 3-bromo-		220	C10H21Br	030571-71-2	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DDCB4DL

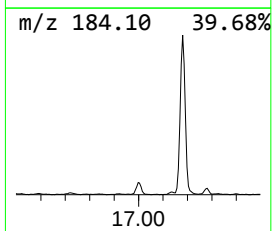
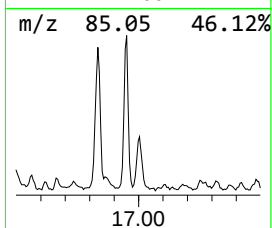
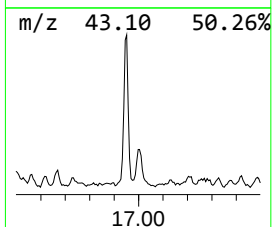
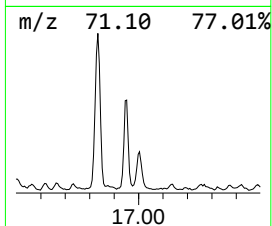
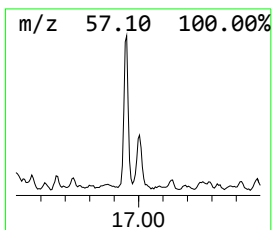
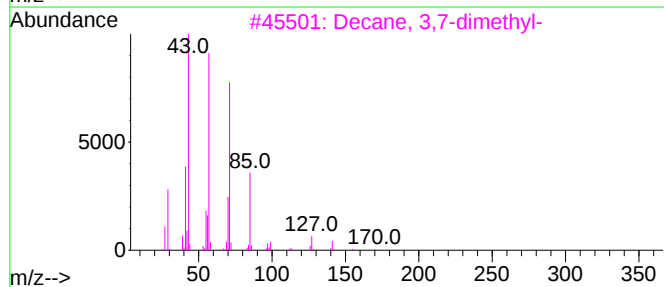
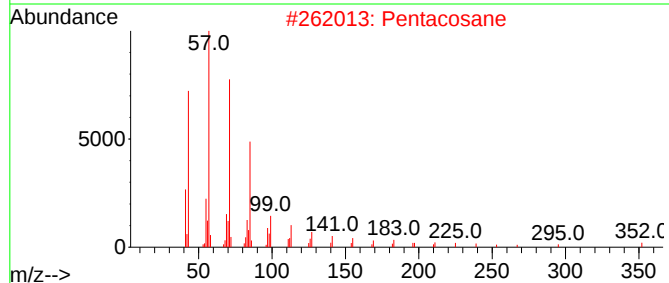
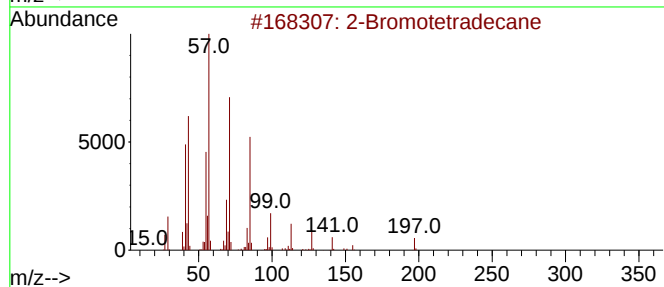
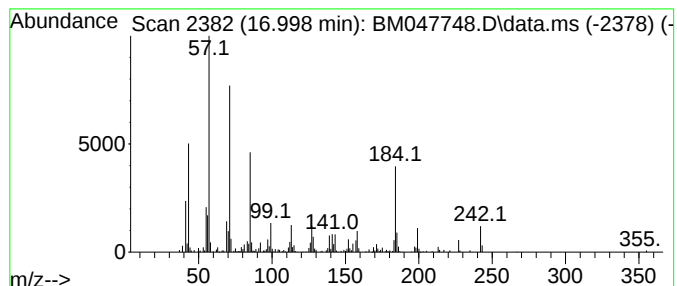
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Quant Title : SVOA CALIBRATION

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

Peak Number 21 2-Bromotetradecane Concentration Rank 21

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.998	2.14 ng/ul	313894	Phenanthrene-d10	17.180

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Bromotetradecane	276	C14H29Br	074036-95-6	62
2			Pentacosane	352	C25H52	000629-99-2	58
3			Decane, 3,7-dimethyl-	170	C12H26	017312-54-8	58
4			Tetracosane, 11-decyl-	479	C34H70	055429-84-0	53
5			Hentriacontane	437	C31H64	000630-04-6	53



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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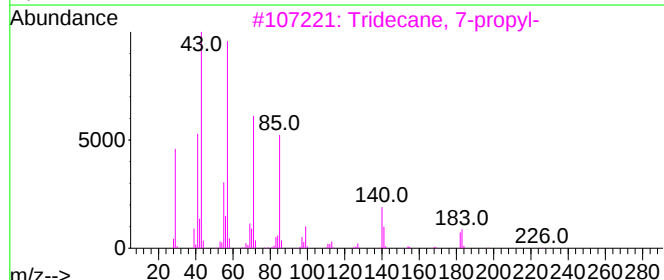
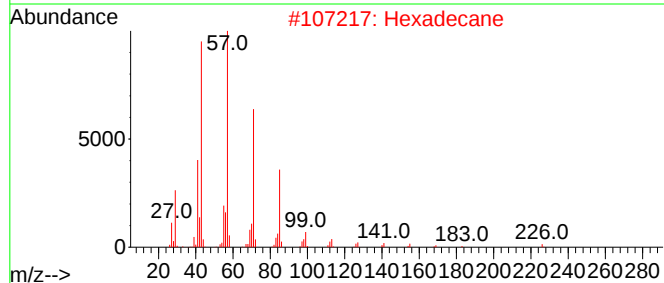
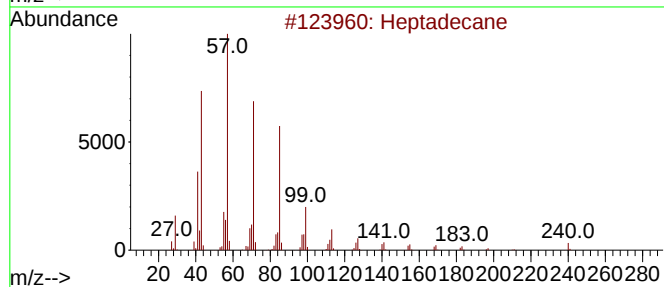
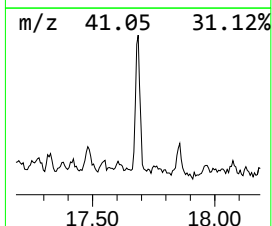
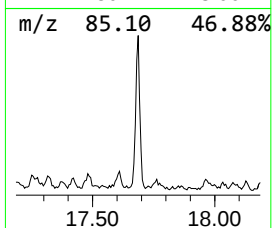
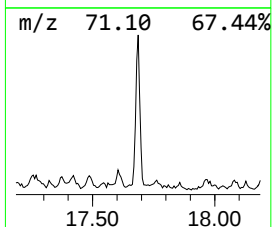
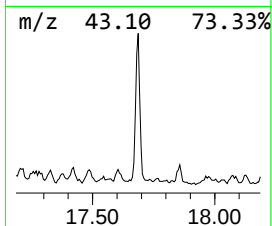
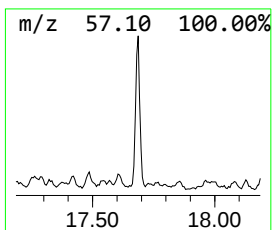
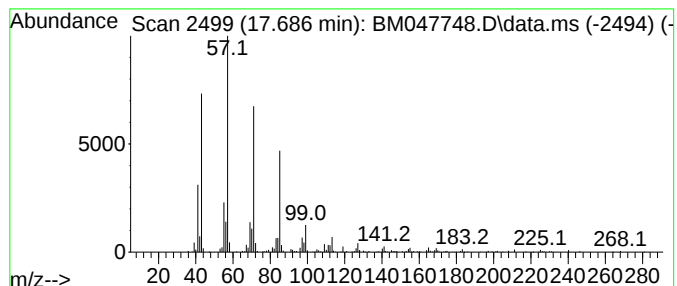
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Quant Title : SVOA CALIBRATION

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

Peak Number 22 (DEL) Alkane: Straight-Chai... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.686	2.20 ng/ul	323463	Phenanthrene-d10	17.180

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	96
2		Hexadecane	226	C16H34	000544-76-3	94
3		Tridecane, 7-propyl-	226	C16H34	055045-09-5	91
4		Nonadecane	268	C19H40	000629-92-5	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM100124\
Data File : BM047748.D
Acq On : 01 Oct 2024 18:21
Operator : RC/JU
Sample : P4018-08DL 30X
Misc :
ALS Vial : 12 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM092724.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
4H-Imidazol-4-o...	6.975	3.0	ng/ul	204385	1	7.810	1353310	20.0
(DEL) Alkane: S...	7.446	9.4	ng/ul	635597	1	7.810	1353310	20.0
1-Hexanol, 2-et...	7.881	2.2	ng/ul	145885	1	7.810	1353310	20.0
(DEL) Alkane: S...	8.351	2.4	ng/ul	163555	1	7.810	1353310	20.0
(DEL) Alkane: S...	8.581	2.9	ng/ul	194165	1	7.810	1353310	20.0
Benzene, 4-ethy...	8.922	3.0	ng/ul	199492	1	7.810	1353310	20.0
(DEL) Alkane: S...	9.046	13.6	ng/ul	919055	1	7.810	1353310	20.0
unknown-01	9.169	3.0	ng/ul	203812	1	7.810	1353310	20.0
4-Nonanone, 2,6...	10.728	3.3	ng/ul	521406	2	10.604	3203400	20.0
unknown-02	10.987	2.8	ng/ul	454061	2	10.604	3203400	20.0
Benzene, 1,2,4-...	11.704	3.5	ng/ul	566992	2	10.604	3203400	20.0
(DEL) Alkane: S...	12.034	3.9	ng/ul	619434	2	10.604	3203400	20.0
unknown-03	12.681	2.3	ng/ul	279283	3	14.445	2388410	20.0
(DEL) Alkane: S...	13.281	5.0	ng/ul	594922	3	14.445	2388410	20.0
Oxalic acid, bi...	14.351	22.8	ng/ul	2718650	3	14.445	2388410	20.0
9-Methyl-S-octa...	14.586	3.6	ng/ul	425337	3	14.445	2388410	20.0
Naphthalene, 2,...	15.204	2.0	ng/ul	242046	3	14.445	2388410	20.0
(DEL) Alkane: S...	15.298	6.1	ng/ul	733461	3	14.445	2388410	20.0
(DEL) Alkane: S...	16.163	4.0	ng/ul	583371	4	17.180	2934280	20.0
(DEL) Alkane: S...	16.951	3.3	ng/ul	488161	4	17.180	2934280	20.0
2-Bromotetradecane	16.998	2.1	ng/ul	313894	4	17.180	2934280	20.0
(DEL) Alkane: S...	17.686	2.2	ng/ul	323463	4	17.180	2934280	20.0