

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
 Data File : BM037848.D
 Acq On : 05 Dec 2022 17:30
 Operator : CG/JU
 Sample : N5738-01DL 5X
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GBNH2DL

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-BM120122.M
 Title : SVOA CALIBRATION

Signal : TIC: BM037848.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.422	4	6	10	rVB3	12113	13797	0.32%	0.038%
2	3.734	52	59	67	rBV5	10817	24543	0.56%	0.068%
3	3.887	81	85	97	rBV	56648	132081	3.02%	0.363%
4	3.975	97	100	103	rVB3	11724	18005	0.41%	0.050%
5	4.105	118	122	126	rVB2	14485	24049	0.55%	0.066%
6	4.199	134	138	146	rVB	21938	33332	0.76%	0.092%
7	4.581	200	203	208	rBV7	4356	7405	0.17%	0.020%
8	5.158	298	301	304	rVB4	6441	7639	0.17%	0.021%
9	5.252	314	317	321	rVB4	4646	5929	0.14%	0.016%
10	5.399	341	342	349	rVB6	5908	8843	0.20%	0.024%
11	5.752	397	402	406	rBV8	6821	10927	0.25%	0.030%
12	6.046	449	452	458	rBV8	5657	13011	0.30%	0.036%
13	6.587	541	544	548	rBV6	7784	10268	0.24%	0.028%
14	6.758	567	573	579	rVB9	6708	13007	0.30%	0.036%
15	6.946	602	605	607	rVB3	4837	5650	0.13%	0.016%
16	7.069	620	626	631	rBV9	9196	19198	0.44%	0.053%
17	7.246	646	656	666	rVB	253115	423326	9.69%	1.164%
18	7.416	672	685	693	rBV2	209573	408823	9.36%	1.124%
19	7.622	712	720	727	rBV	236644	382703	8.76%	1.053%
20	8.093	792	800	807	rBV	1177662	1781262	40.77%	4.899%
21	8.152	807	810	818	rVB3	11764	19162	0.44%	0.053%
22	8.399	851	852	858	rVB6	3571	5340	0.12%	0.015%
23	8.763	910	914	915	rBV4	5313	6502	0.15%	0.018%
24	8.804	915	921	927	rVB	226453	360359	8.25%	0.991%
25	8.922	936	941	948	rVB	58334	93467	2.14%	0.257%
26	9.034	957	960	961	rBV3	5621	6189	0.14%	0.017%
27	9.263	992	999	1007	rVB	242004	388172	8.88%	1.068%
28	9.316	1007	1008	1012	rBV4	4378	5865	0.13%	0.016%
29	9.363	1012	1016	1021	rVB7	7300	12255	0.28%	0.034%
30	9.657	1065	1066	1068	rBV2	6187	5283	0.12%	0.015%
31	9.693	1070	1072	1076	rVB5	6566	9342	0.21%	0.026%
32	9.810	1088	1092	1096	rBV7	6069	9854	0.23%	0.027%
33	9.987	1115	1122	1130	rBV	179487	308705	7.07%	0.849%
34	10.134	1141	1147	1152	rBV4	17498	32846	0.75%	0.090%
35	10.204	1156	1159	1165	rBV8	6989	12033	0.28%	0.033%
36	10.351	1181	1184	1189	rVB7	5593	7772	0.18%	0.021%

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Integrator: RTE

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Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
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37	10.528	1208	1214	1224	rBV	268232	456140	10.44%	1.255%
38	10.681	1234	1240	1247	rBV	49980	87016	1.99%	0.239%
39	10.851	1264	1269	1270	rBV5	3981	6038	0.14%	0.017%
40	10.916	1270	1280	1287	rBV	1661796	2679612	61.33%	7.370%
41	11.051	1297	1303	1312	rBV2	51711	105948	2.42%	0.291%
42	11.363	1354	1356	1360	rBV4	5260	6274	0.14%	0.017%
43	11.404	1362	1363	1367	rVV3	4958	7031	0.16%	0.019%
44	11.445	1367	1370	1376	rVB8	9941	11954	0.27%	0.033%
45	11.587	1392	1394	1401	rVB8	4167	7435	0.17%	0.020%
46	11.692	1406	1412	1417	rBV10	11919	26211	0.60%	0.072%
47	11.792	1426	1429	1431	rBV3	4539	6266	0.14%	0.017%
48	12.204	1495	1499	1502	rBV6	3306	5647	0.13%	0.016%
49	12.269	1506	1510	1511	rVV4	5257	7119	0.16%	0.020%
50	12.316	1514	1518	1520	rVV5	5172	6630	0.15%	0.018%
51	12.492	1546	1548	1554	rVB7	8036	10865	0.25%	0.030%
52	12.563	1557	1560	1565	rBV7	5505	6456	0.15%	0.018%
53	12.645	1572	1574	1576	rBV3	4049	5303	0.12%	0.015%
54	12.881	1610	1614	1618	rBV7	6617	11280	0.26%	0.031%
55	12.957	1622	1627	1631	rVV2	30911	45875	1.05%	0.126%
56	13.092	1644	1650	1654	rVB7	18107	27911	0.64%	0.077%
57	13.157	1658	1661	1666	rVV7	6163	12412	0.28%	0.034%
58	13.198	1666	1668	1670	rVV3	5022	6183	0.14%	0.017%
59	13.239	1671	1675	1680	rVB7	18189	27968	0.64%	0.077%
60	13.486	1713	1717	1720	rBV6	9278	15571	0.36%	0.043%
61	13.534	1723	1725	1728	rVB4	8820	8146	0.19%	0.022%
62	13.604	1732	1737	1742	rBV5	25681	48027	1.10%	0.132%
63	13.734	1755	1759	1764	rBV2	37261	50881	1.16%	0.140%
64	13.863	1777	1781	1784	rBV6	15983	28273	0.65%	0.078%
65	13.910	1784	1789	1796	rVB10	23839	47012	1.08%	0.129%
66	13.998	1797	1804	1808	rBV	459306	673498	15.41%	1.852%
67	14.039	1808	1811	1819	rVB2	125783	186782	4.27%	0.514%
68	14.128	1819	1826	1834	rBV	452189	686228	15.71%	1.887%
69	14.245	1841	1846	1851	rVB2	85578	117505	2.69%	0.323%
70	14.416	1869	1875	1884	rBV	572622	898385	20.56%	2.471%
71	14.557	1892	1899	1903	rBV3	63892	149056	3.41%	0.410%
72	14.604	1903	1907	1915	rVB	174508	257056	5.88%	0.707%
73	14.722	1921	1927	1935	rBV	2334243	3397650	77.76%	9.345%
74	14.916	1955	1960	1965	rBV	122735	202030	4.62%	0.556%
75	14.963	1965	1968	1974	rVB	90494	126098	2.89%	0.347%
76	15.069	1984	1986	1989	rBV4	11833	14971	0.34%	0.041%

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77	15.304	2023	2026	2028	rBV3	18604	23605	0.54%	0.065%
78	15.339	2028	2032	2037	rVB4	43063	66251	1.52%	0.182%
79	15.545	2062	2067	2071	rVB3	43573	76709	1.76%	0.211%
80	15.710	2089	2095	2106	rBV	694696	1048591	24.00%	2.884%
81	15.822	2109	2114	2120	rVB2	124621	193852	4.44%	0.533%
82	15.945	2132	2135	2139	rVV3	38133	49339	1.13%	0.136%
83	15.998	2139	2144	2154	rVB	345320	469745	10.75%	1.292%
84	16.404	2210	2213	2215	rVB2	78234	99250	2.27%	0.273%
85	17.110	2327	2333	2343	rBV	3006747	4369351	100.00%	12.017%
86	17.198	2345	2348	2351	rVB	79089	90612	2.07%	0.249%
87	17.339	2369	2372	2376	rVB	80081	96528	2.21%	0.265%
88	17.463	2387	2393	2401	rBV	2572807	3783288	86.59%	10.405%
89	17.557	2405	2409	2420	rVB	666440	985264	22.55%	2.710%
90	17.933	2471	2473	2478	rVB2	95610	107079	2.45%	0.295%
91	18.221	2519	2522	2526	rBV2	127324	158068	3.62%	0.435%
92	18.339	2538	2542	2547	rBV	619607	836266	19.14%	2.300%
93	18.627	2588	2591	2597	rVB3	136911	196591	4.50%	0.541%
94	19.604	2754	2757	2761	rBV2	228520	278902	6.38%	0.767%
95	19.839	2793	2797	2805	rVB	752105	1018639	23.31%	2.802%
96	20.704	2941	2944	2948	rVV	771301	861238	19.71%	2.369%
97	20.745	2948	2951	2957	rVB	505185	630997	14.44%	1.735%
98	21.309	3044	3047	3051	rBV	398330	471825	10.80%	1.298%
99	21.627	3097	3101	3107	rBV	2093871	2664111	60.97%	7.327%
100	24.115	3519	3524	3531	rVB2	1362163	2674967	61.22%	7.357%

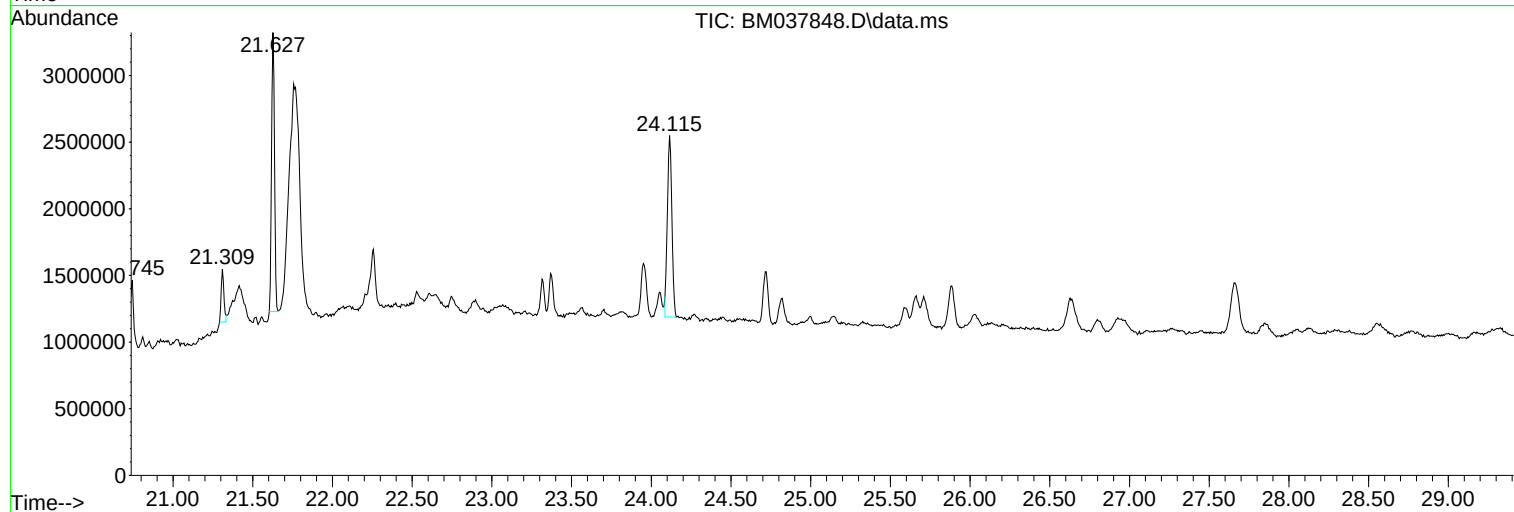
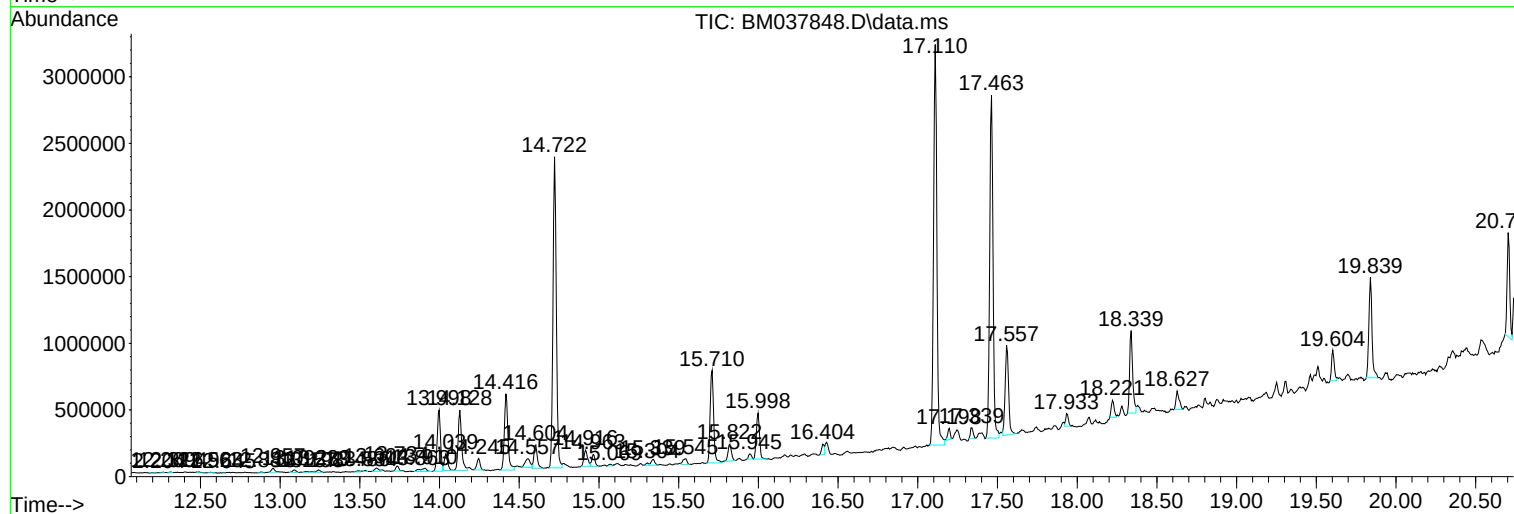
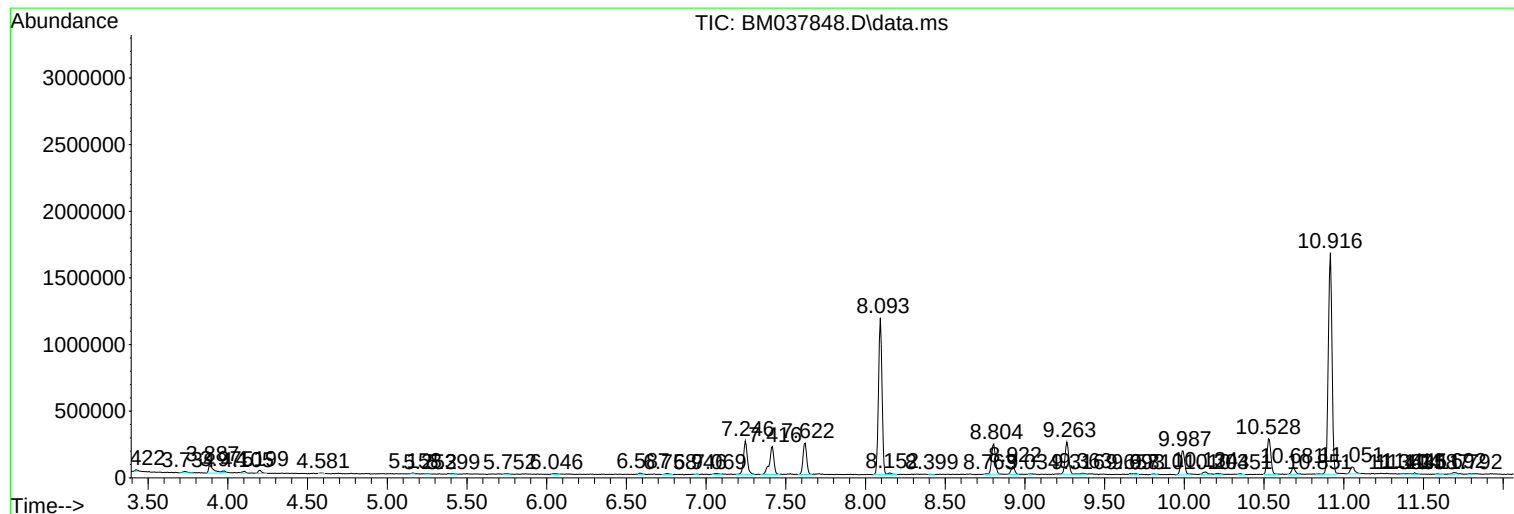
Sum of corrected areas: 36358755

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ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNH2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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



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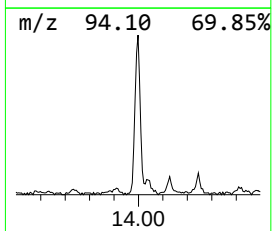
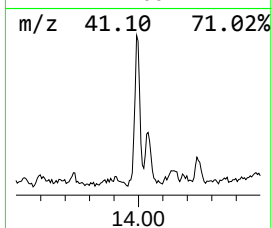
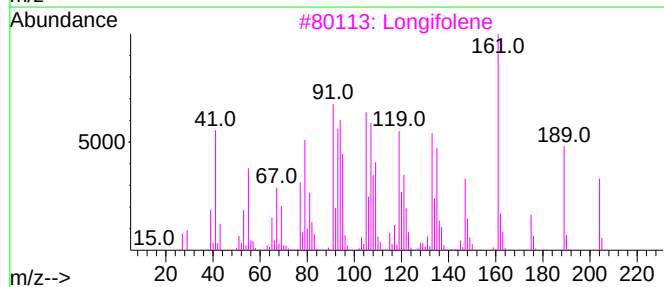
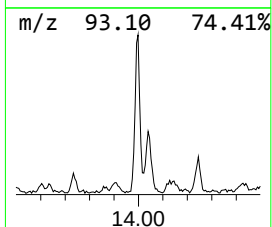
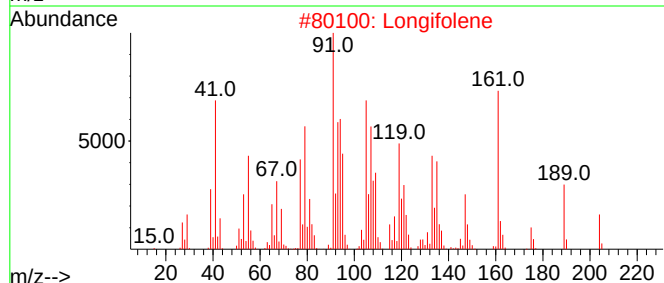
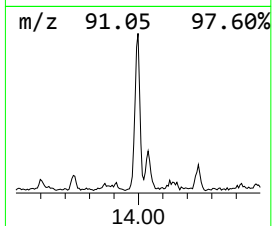
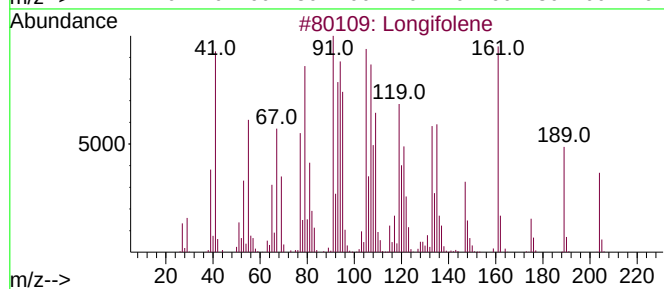
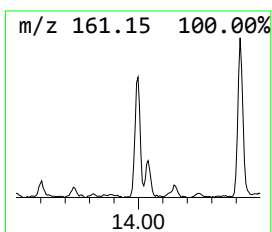
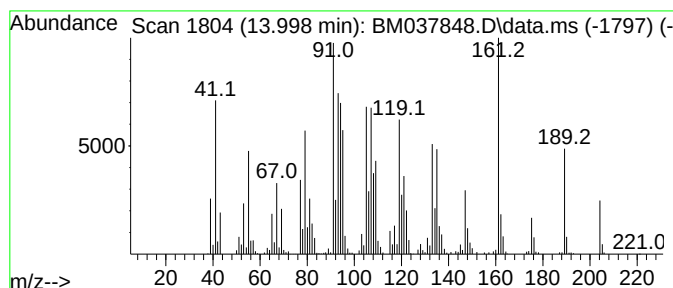
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Longifolene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.998	3.96 ng/ul	673498	Acenaphthene-d10	14.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Longifolene	204	C15H24	000475-20-7	99
2			Longifolene	204	C15H24	000475-20-7	99
3			Longifolene	204	C15H24	000475-20-7	99
4			Longifolene	204	C15H24	000475-20-7	99
5			Longifolene	204	C15H24	000475-20-7	96



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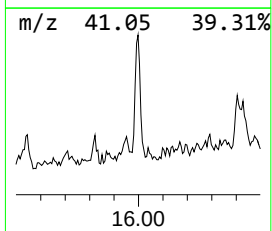
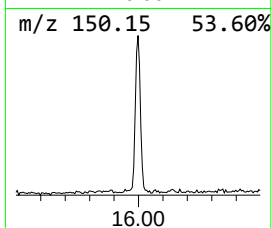
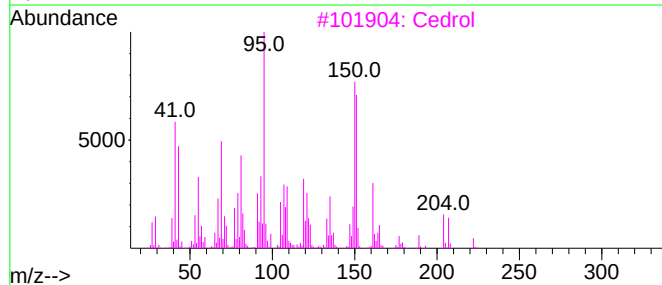
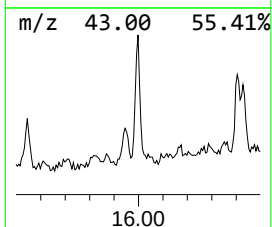
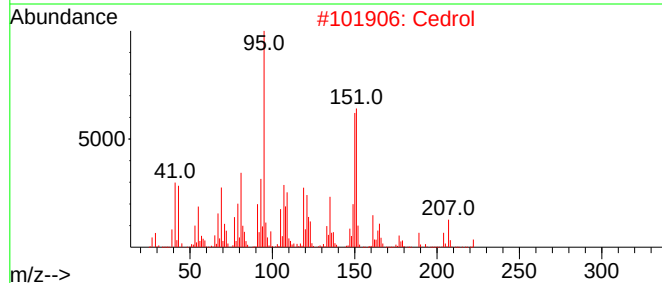
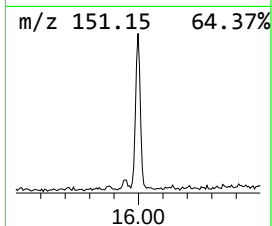
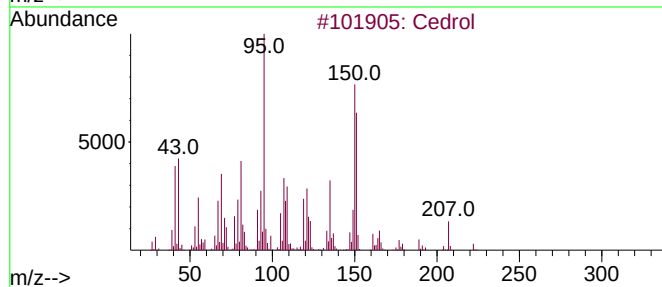
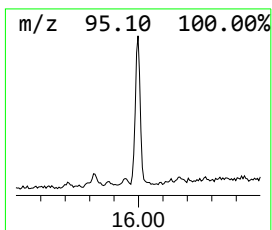
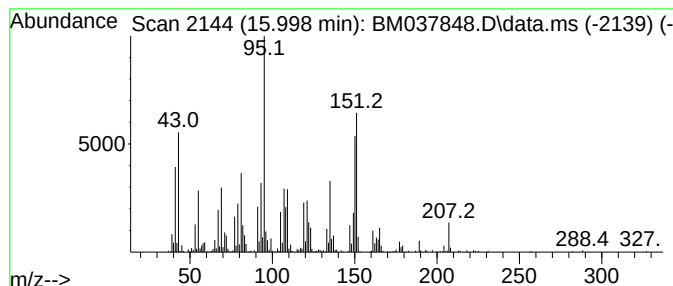
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cedrol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.998	2.77 ng/ul	469745	Acenaphthene-d10	14.722

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cedrol	222	C15H26O	000077-53-2	94
2			Cedrol	222	C15H26O	000077-53-2	91
3			Cedrol	222	C15H26O	000077-53-2	87
4			(3R,3aS,6S,7R)-3,6,8,8-Tetrameth...	222	C15H26O	019903-73-2	87
5			(3R,3aS,6S,7R)-3,6,8,8-Tetrameth...	222	C15H26O	019903-73-2	72



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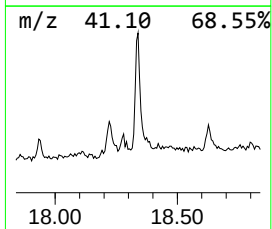
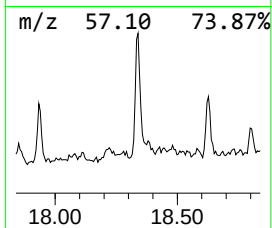
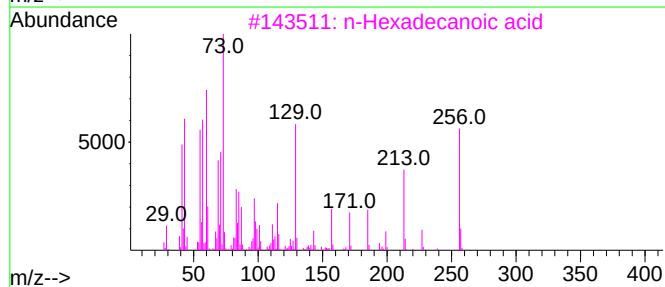
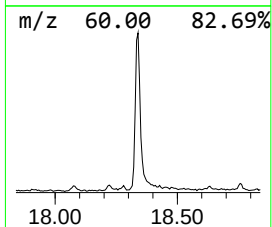
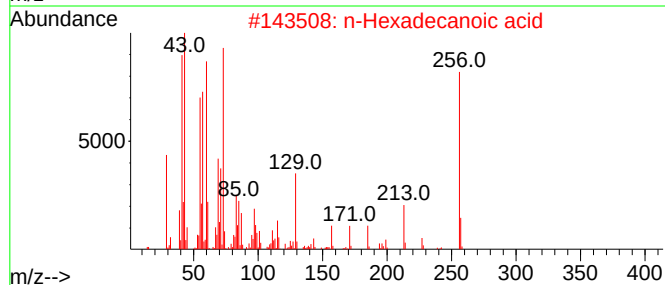
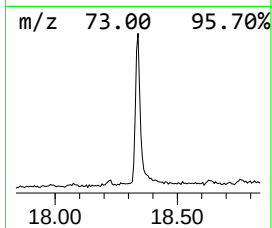
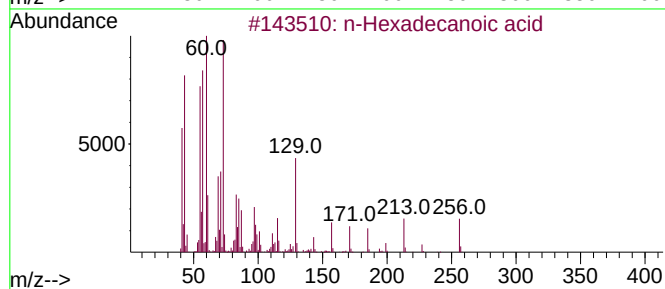
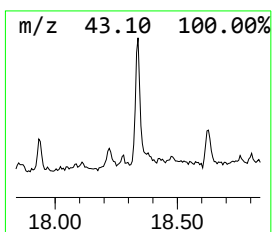
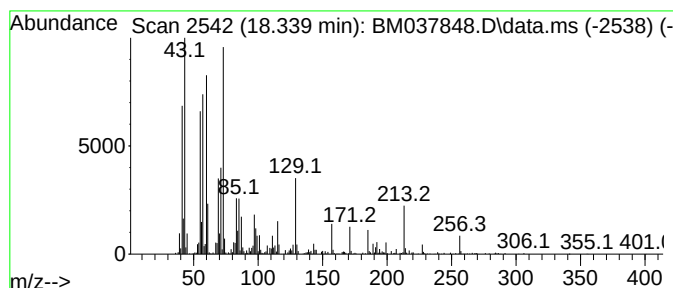
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
18.339	4.42 ng/ul	836266	Phenanthrene-d10		17.463	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	99
2	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	97
3	n-Hexadecanoic acid		256	C16H32O2	000057-10-3	96
4	Tridecanoic acid		214	C13H26O2	000638-53-9	95
5	Tridecanoic acid		214	C13H26O2	000638-53-9	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
Data File : BM037848.D
Acq On : 05 Dec 2022 17:30
Operator : CG/JU
Sample : N5738-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNH2DL

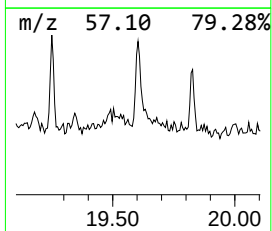
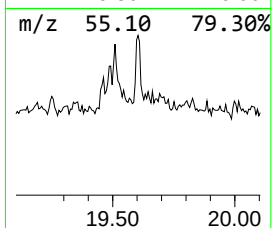
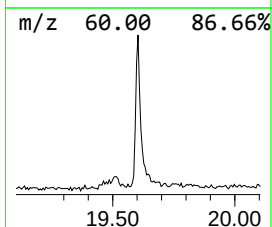
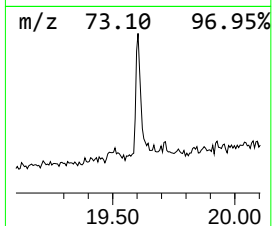
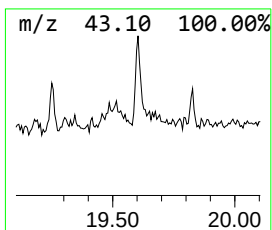
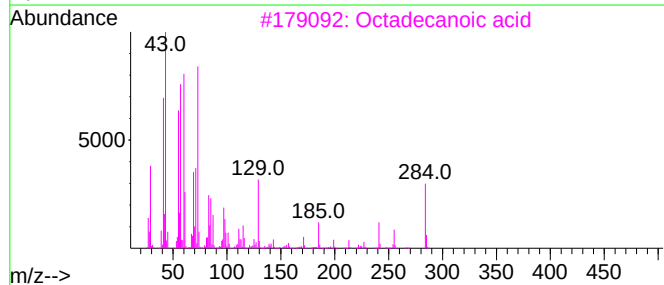
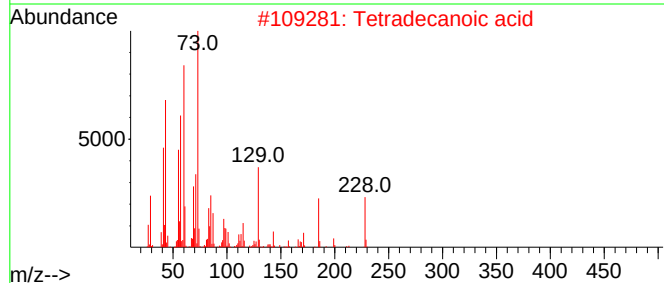
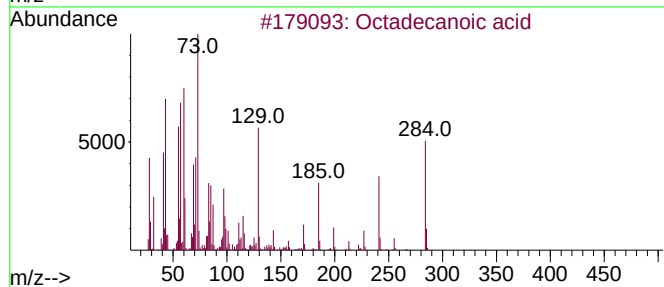
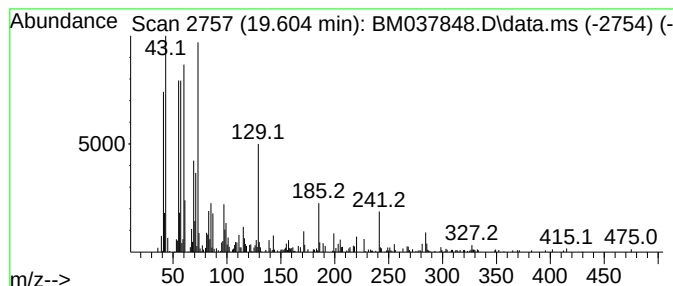
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.604	2.09 ng/ul	278902	Chrysene-d12	21.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	97
2			Tetradecanoic acid	228	C14H28O2	000544-63-8	87
3			Octadecanoic acid	284	C18H36O2	000057-11-4	80
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	76
5			Tridecanoic acid	214	C13H26O2	000638-53-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
Data File : BM037848.D
Acq On : 05 Dec 2022 17:30
Operator : CG/JU
Sample : N5738-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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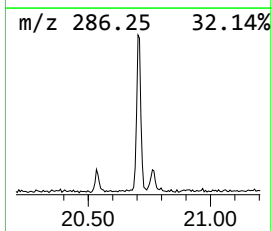
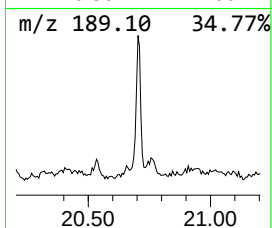
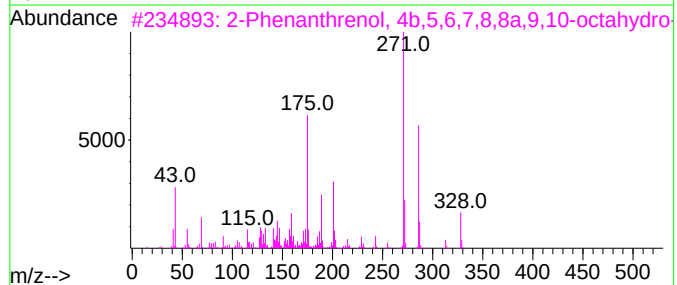
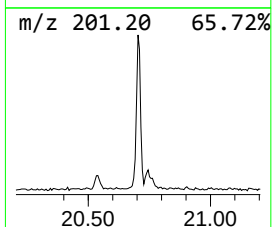
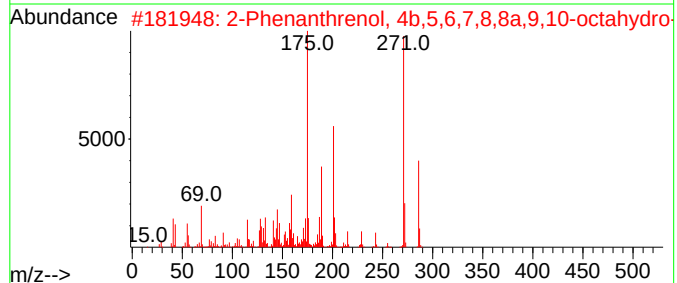
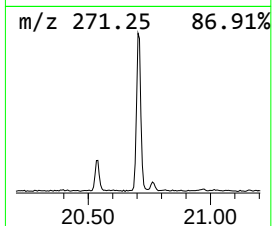
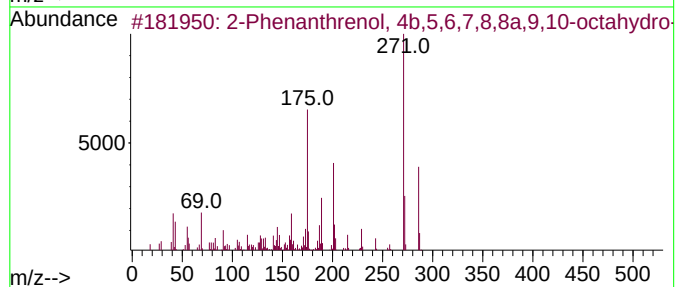
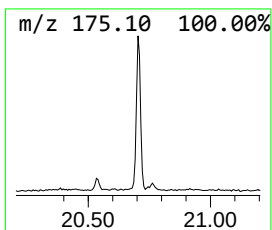
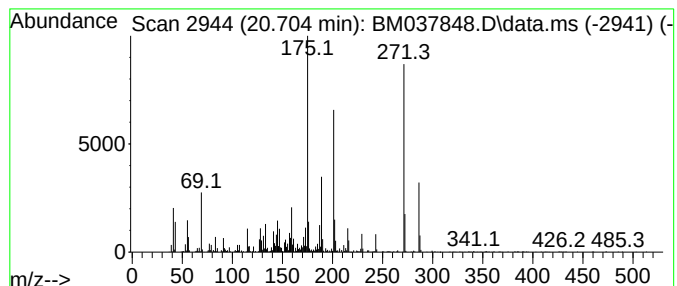
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 2-Phenanthrenol, 4b,5,6,7,8... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.704	6.47 ng/ul	861238	Chrysene-d12	21.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	96		
2	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	286	C20H30O	000511-15-9	95		
3	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	91		
4	2-Phenanthrenol, 4b,5,6,7,8,8a,9...	328	C22H32O2	015340-82-6	46		
5	Pisiferol	302	C20H30O2	024035-36-7	46		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
Data File : BM037848.D
Acq On : 05 Dec 2022 17:30
Operator : CG/JU
Sample : N5738-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNH2DL

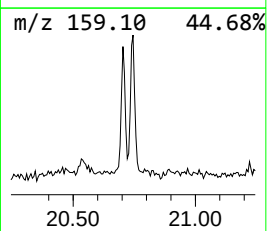
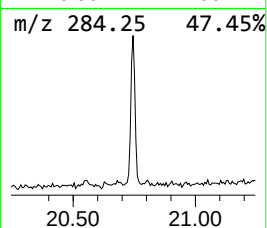
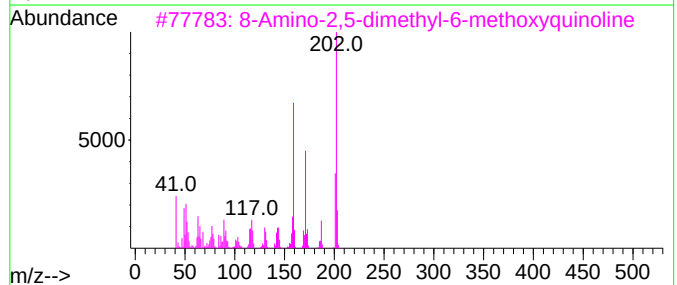
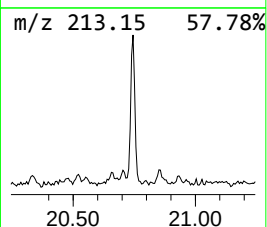
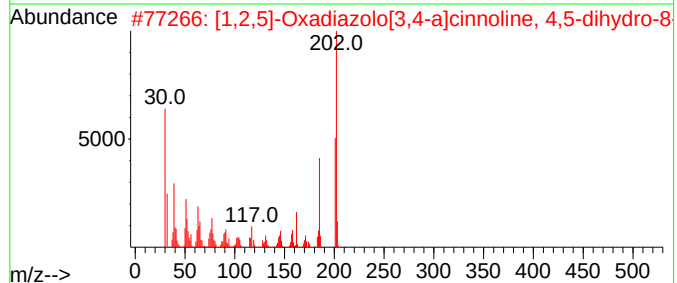
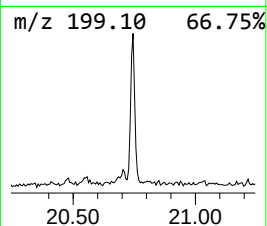
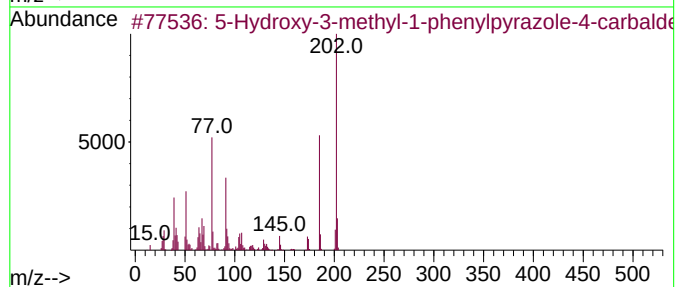
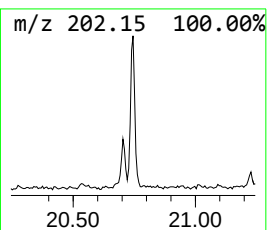
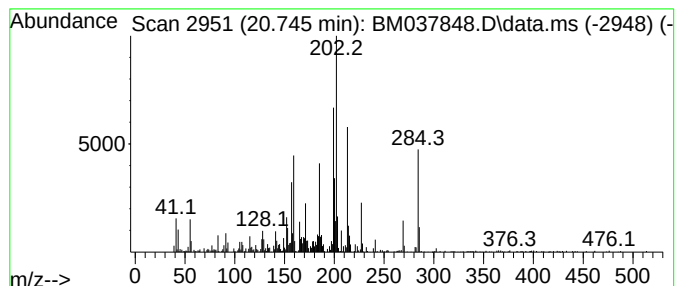
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.745	4.74 ng/ul	630997	Chrysene-d12	21.627

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			5-Hydroxy-3-methyl-1-phenylpyraz...	202	C11H10N2O2	060484-29-9	38
2			[1,2,5]-Oxadiazolo[3,4-a]cinnoli...	202	C10H10N4O	1000260-59-4	30
3			8-Amino-2,5-dimethyl-6-methoxyqu...	202	C12H14N2O	1000214-69-9	25
4			4,4'-Bis(tetrahydrothiopyran)	202	C10H18S2	116196-83-9	25
5			8-Hydroxy-naphthoic acid, methyl...	202	C12H10O3	005991-56-0	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
Data File : BM037848.D
Acq On : 05 Dec 2022 17:30
Operator : CG/JU
Sample : N5738-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

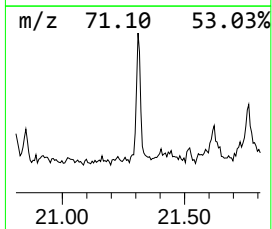
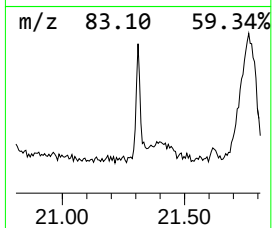
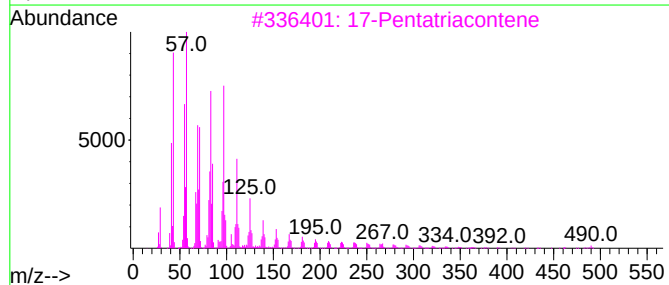
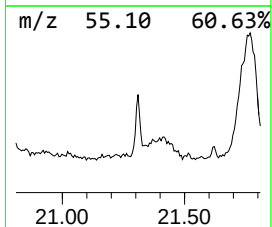
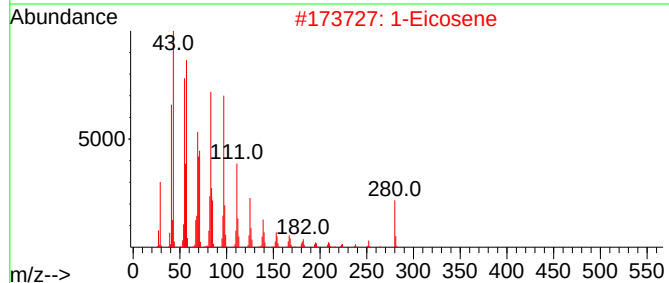
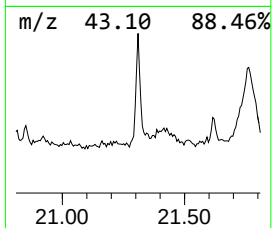
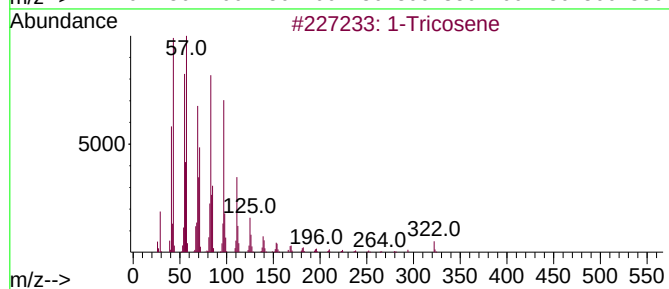
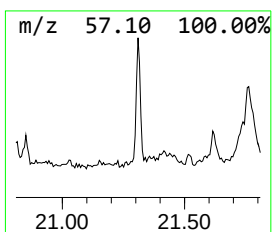
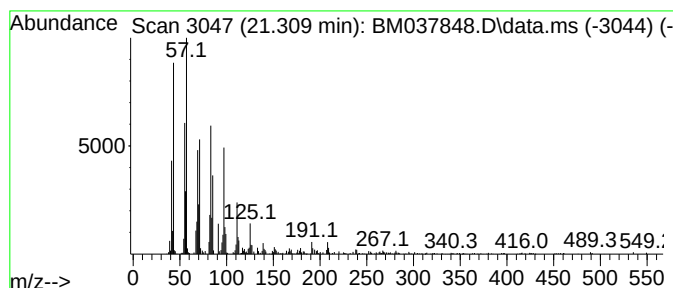
Instrument :
BNA_M
ClientSampleId :
GBNH2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
21.310	3.54 ng/ul	471825	Chrysene-d12		21.627	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Tricosene		322	C23H46	018835-32-0	96
2	1-Eicosene		280	C20H40	003452-07-1	95
3	17-Pentatriacontene		491	C35H70	006971-40-0	94
4	9-Tricosene, (Z)-		322	C23H46	027519-02-4	93
5	9-Tricosene, (Z)-		322	C23H46	027519-02-4	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120522\
Data File : BM037848.D
Acq On : 05 Dec 2022 17:30
Operator : CG/JU
Sample : N5738-01DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNH2DL

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.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
Longifolene	13.998	4.0	ng/ul	673498	3	14.722	3397650	20.0
Cedrol	15.998	2.8	ng/ul	469745	3	14.722	3397650	20.0
n-Hexadecanoic ...	18.339	4.4	ng/ul	836266	4	17.463	3783290	20.0
Octadecanoic acid	19.604	2.1	ng/ul	278902	5	21.627	2664110	20.0
2-Phenanthrenol...	20.704	6.5	ng/ul	861238	5	21.627	2664110	20.0
unknown-01	20.745	4.7	ng/ul	630997	5	21.627	2664110	20.0
(DEL) Alkane: S...	21.310	3.5	ng/ul	471825	5	21.627	2664110	20.0