

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
 Data File : BM037864.D
 Acq On : 06 Dec 2022 21:12
 Operator : CG/JU
 Sample : N5738-12MEDL 2X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 GBNJ1MEDL

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: BM037864.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.428	4	7	11	rVB2	21953	27348	0.41%	0.056%
2	4.098	117	121	125	rVB	21056	28694	0.43%	0.058%
3	4.198	134	138	142	rBV	20051	28160	0.42%	0.057%
4	4.428	175	177	182	rBV6	5114	8489	0.13%	0.017%
5	4.581	202	203	209	rVB5	9361	11296	0.17%	0.023%
6	5.157	294	301	309	rBV5	14341	33801	0.51%	0.069%
7	6.698	559	563	566	rBV5	4534	6753	0.10%	0.014%
8	7.239	649	655	667	rVB	350758	567203	8.50%	1.153%
9	7.410	676	684	696	rVB	312925	485113	7.27%	0.986%
10	7.616	712	719	725	rBV	371810	569248	8.53%	1.157%
11	7.698	729	733	737	rVB7	5883	6713	0.10%	0.014%
12	8.092	793	800	807	rBV	940909	1489743	22.33%	3.028%
13	8.151	807	810	817	rVB8	7038	11131	0.17%	0.023%
14	8.798	909	920	930	rBV	399593	634945	9.52%	1.291%
15	8.922	936	941	948	rVB2	18079	32655	0.49%	0.066%
16	9.263	992	999	1006	rBV	322283	539925	8.09%	1.098%
17	9.657	1062	1066	1074	rVB3	13987	23886	0.36%	0.049%
18	9.845	1096	1098	1104	rBV4	6067	11100	0.17%	0.023%
19	9.986	1115	1122	1132	rBV	239561	395475	5.93%	0.804%
20	10.527	1207	1214	1223	rBV	380698	626987	9.40%	1.275%
21	10.910	1271	1279	1289	rVB	1264983	2049335	30.71%	4.166%
22	11.051	1295	1303	1310	rBV	119802	226781	3.40%	0.461%
23	11.263	1334	1339	1345	rVB9	5317	11135	0.17%	0.023%
24	11.386	1353	1360	1361	rBV7	5572	8785	0.13%	0.018%
25	11.592	1391	1395	1400	rBV7	4686	9274	0.14%	0.019%
26	11.692	1407	1412	1416	rBV8	5144	8906	0.13%	0.018%
27	11.816	1429	1433	1436	rBV6	4787	7580	0.11%	0.015%
28	11.916	1446	1450	1455	rBV3	19535	28122	0.42%	0.057%
29	11.957	1455	1457	1462	rVV5	6643	8943	0.13%	0.018%
30	12.304	1512	1516	1520	rBV2	45186	63236	0.95%	0.129%
31	12.410	1529	1534	1537	rBV7	8023	14770	0.22%	0.030%
32	12.492	1543	1548	1551	rBV4	7710	14720	0.22%	0.030%
33	12.633	1566	1572	1574	rBV6	5105	7434	0.11%	0.015%
34	12.663	1574	1577	1581	rVB6	5816	8105	0.12%	0.016%
35	12.886	1609	1615	1618	rBV8	4222	7199	0.11%	0.015%
36	12.933	1618	1623	1626	rBV6	15647	24214	0.36%	0.049%

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Integrator: RTE

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Stop Thrs : 0

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

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

37	13.039	1637	1641	1644	rVB4	10151	14478	0.22%	0.029%
38	13.110	1649	1653	1658	rBV4	14040	21270	0.32%	0.043%
39	13.192	1665	1667	1671	rVB4	11669	12313	0.18%	0.025%
40	13.245	1672	1676	1682	rVB	55421	76603	1.15%	0.156%
41	13.533	1716	1725	1733	rBV	128379	210543	3.16%	0.428%
42	13.610	1733	1738	1745	rBV5	25437	57778	0.87%	0.117%
43	13.745	1757	1761	1762	rBV4	5730	6995	0.10%	0.014%
44	13.774	1762	1766	1769	rVB6	10454	15696	0.24%	0.032%
45	13.921	1788	1791	1795	rVB6	9441	13954	0.21%	0.028%
46	14.045	1809	1812	1814	rBV4	12661	18454	0.28%	0.038%
47	14.121	1818	1825	1832	rVV	544207	852268	12.77%	1.733%
48	14.186	1832	1836	1840	rVV	137554	195343	2.93%	0.397%
49	14.227	1840	1843	1849	rVB4	30857	56053	0.84%	0.114%
50	14.304	1854	1856	1860	rVB2	25066	26265	0.39%	0.053%
51	14.357	1862	1865	1868	rBV5	14812	20764	0.31%	0.042%
52	14.415	1869	1875	1880	rBV	731942	1085564	16.27%	2.207%
53	14.545	1893	1897	1901	rVB4	41356	60348	0.90%	0.123%
54	14.598	1901	1906	1914	rBV	951816	1175906	17.62%	2.390%
55	14.668	1914	1918	1921	rBV5	39347	66258	0.99%	0.135%
56	14.721	1921	1927	1938	rVB	1637180	2451764	36.75%	4.984%
57	14.910	1955	1959	1971	rVB2	155612	312780	4.69%	0.636%
58	15.092	1988	1990	1996	rVB3	56056	81446	1.22%	0.166%
59	15.209	2006	2010	2016	rBV3	100434	153164	2.30%	0.311%
60	15.333	2028	2031	2035	rVB4	50103	70357	1.05%	0.143%
61	15.539	2060	2066	2071	rVB4	409978	826663	12.39%	1.680%
62	15.645	2081	2084	2088	rBV4	47565	68623	1.03%	0.140%
63	15.704	2089	2094	2101	rBV	947525	1398149	20.95%	2.842%
64	15.815	2109	2113	2122	rVB3	230265	385142	5.77%	0.783%
65	15.945	2129	2135	2140	rBV	433884	710747	10.65%	1.445%
66	16.098	2158	2161	2168	rBV5	68429	146998	2.20%	0.299%
67	16.162	2169	2172	2176	rVB3	103698	127736	1.91%	0.260%
68	16.404	2209	2213	2214	rBV	682639	766055	11.48%	1.557%
69	16.427	2214	2217	2222	rVB	1136971	1515030	22.71%	3.080%
70	17.109	2327	2333	2343	rBV	4479644	6672348	100.00%	13.564%
71	17.192	2343	2347	2351	rVV	707012	849877	12.74%	1.728%
72	17.245	2351	2356	2361	rVB	1001931	1412358	21.17%	2.871%
73	17.456	2386	2392	2398	rBV	2103980	3119757	46.76%	6.342%
74	17.556	2404	2409	2416	rVB	1195903	1724614	25.85%	3.506%
75	17.850	2456	2459	2465	rVB	291598	420180	6.30%	0.854%
76	17.933	2469	2473	2479	rVB	777188	928322	13.91%	1.887%

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 Title : SVOA CALIBRATION

77	18.333	2537	2541	2545	rBV	234712	341625	5.12%	0.694%
78	18.621	2586	2590	2595	rVB	605721	701490	10.51%	1.426%
79	19.245	2693	2696	2700	rVB	429125	513590	7.70%	1.044%
80	19.833	2791	2796	2808	rVB	1512383	2347626	35.18%	4.772%
81	20.350	2882	2884	2890	rVB2	169273	188141	2.82%	0.382%
82	21.303	3043	3046	3053	rBV2	402870	529607	7.94%	1.077%
83	21.621	3095	3100	3106	rVB	2307791	2981230	44.68%	6.060%
84	23.944	3489	3495	3504	rVB3	1004196	2220969	33.29%	4.515%
85	24.103	3516	3522	3532	rVB2	1547583	3201349	47.98%	6.508%

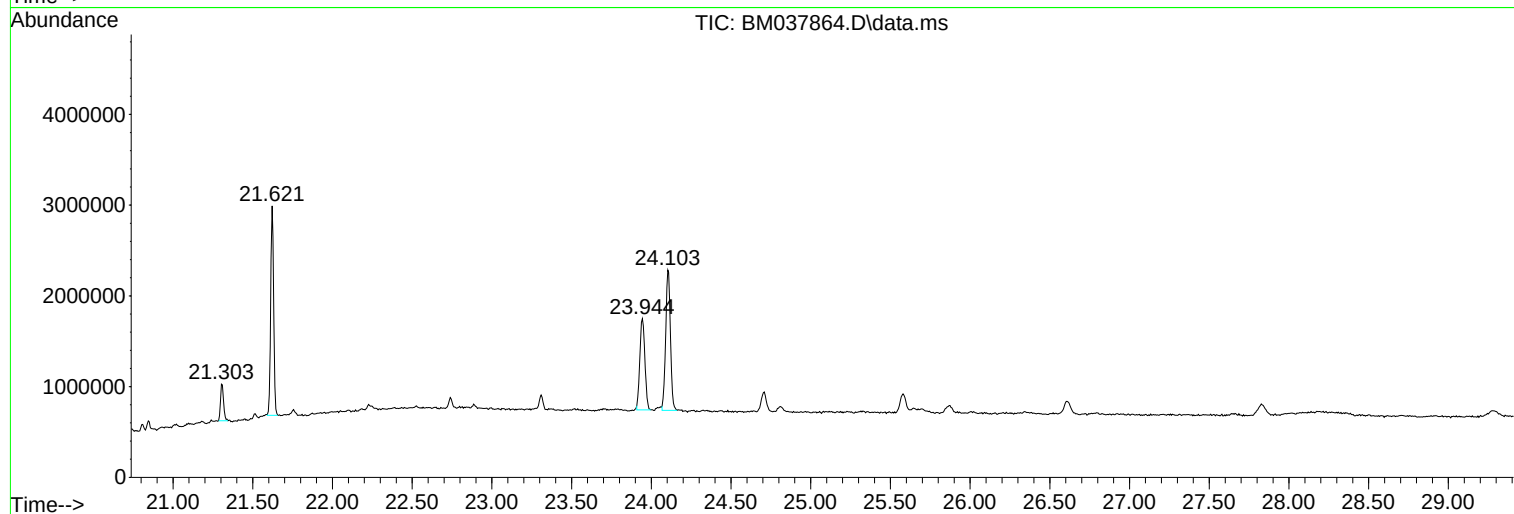
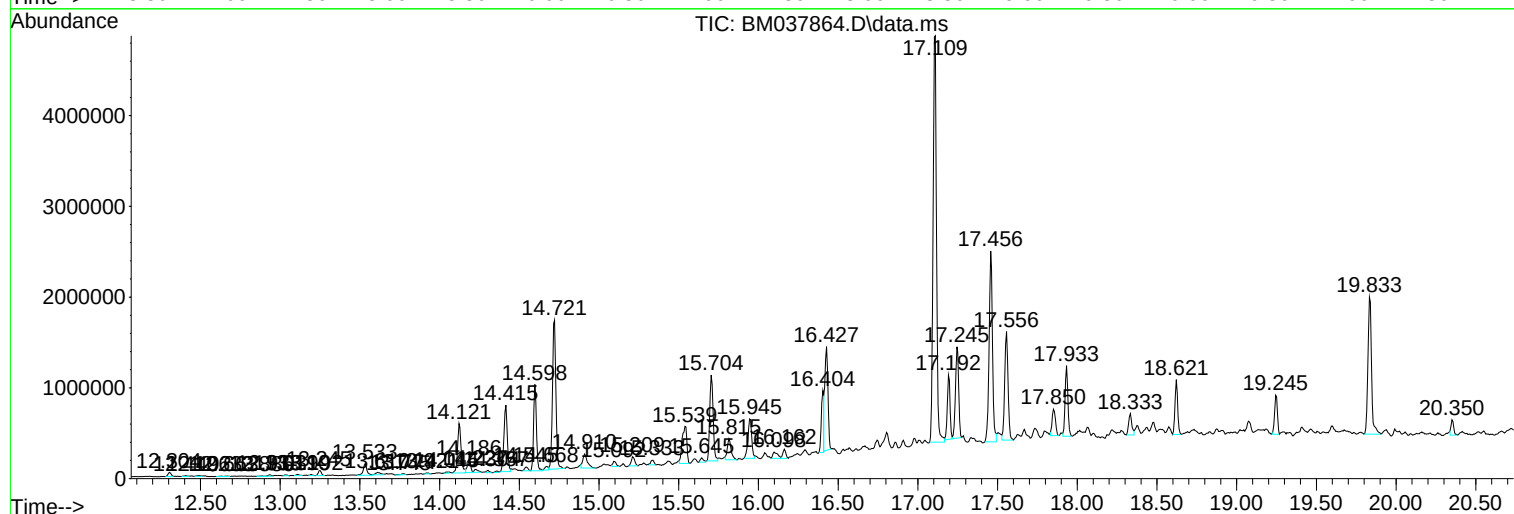
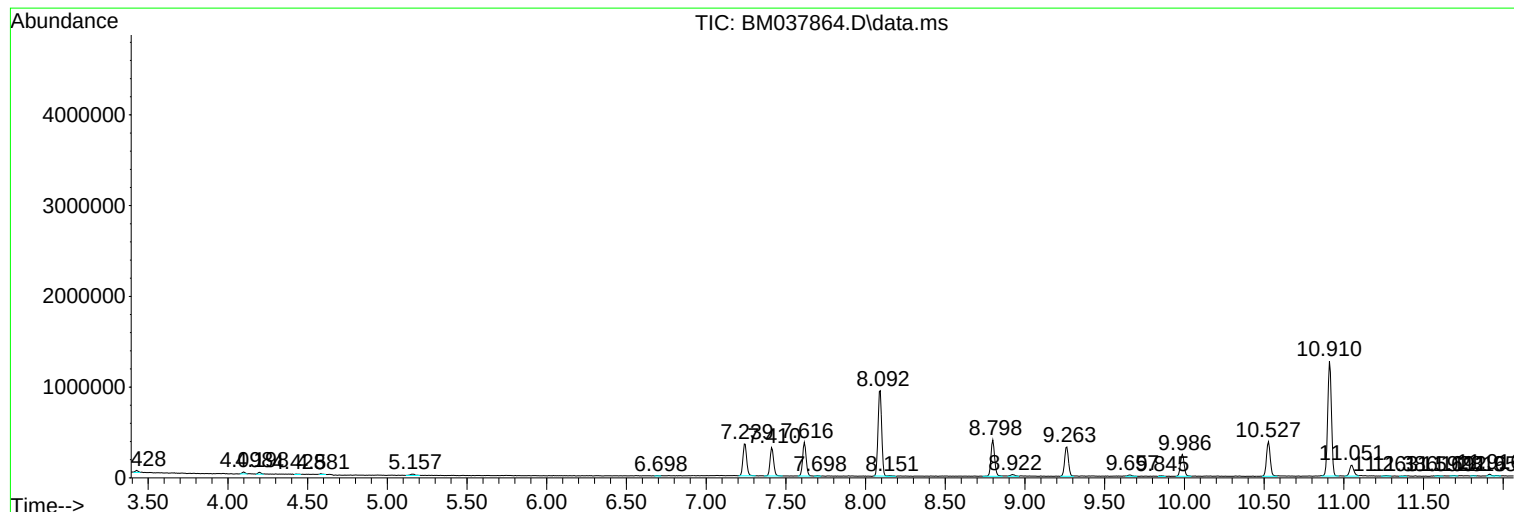
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ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
GBNJ1MEDL

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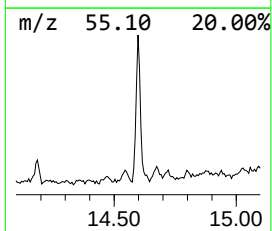
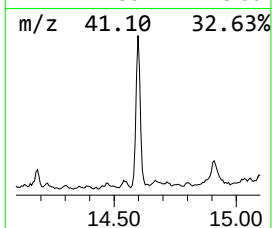
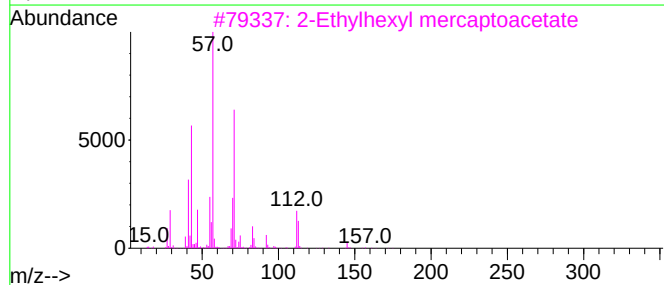
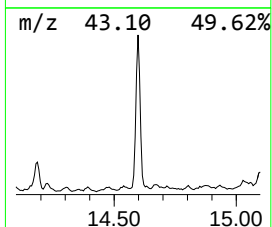
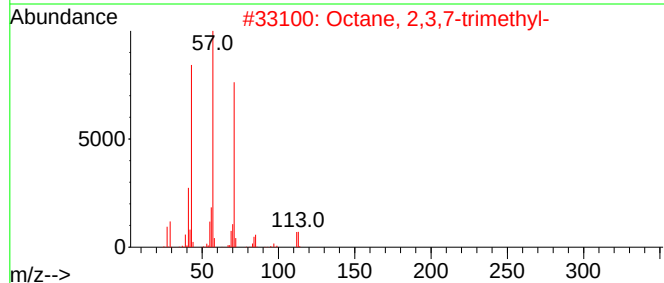
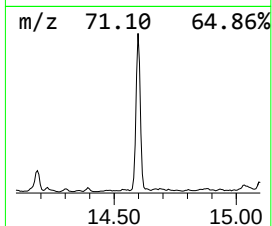
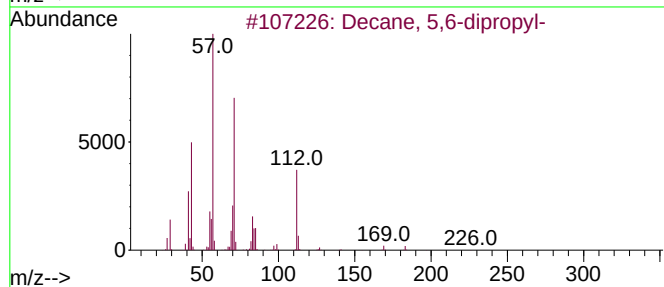
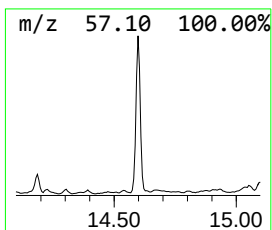
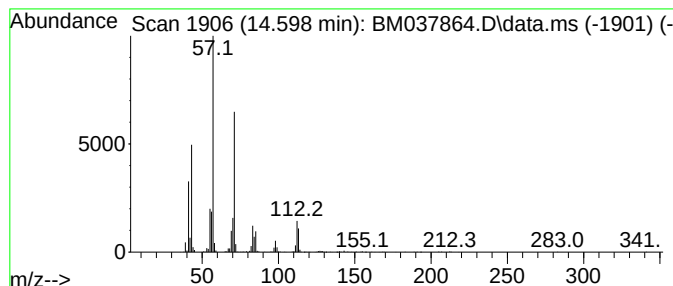
Instrument :
BNA_M
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
14.598	9.59 ng/ul	1175910	Acenaphthene-d10	14.721	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Decane, 5,6-dipropyl-	226	C16H34	119209-20-0	72
2	Octane, 2,3,7-trimethyl-	156	C11H24	062016-34-6	64
3	2-Ethylhexyl mercaptoacetate	204	C10H20O2S	007659-86-1	64
4	Heptane, 3-ethyl-5-methyl-	142	C10H22	052896-90-9	59
5	Carbonic acid, neopentyl 2-ethyl...	244	C14H28O3	1000357-85-7	59



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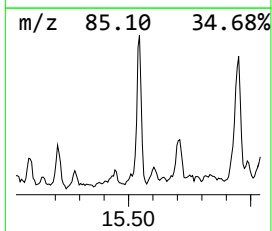
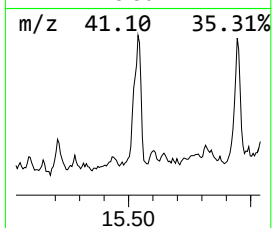
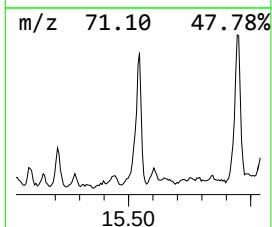
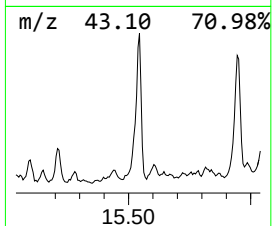
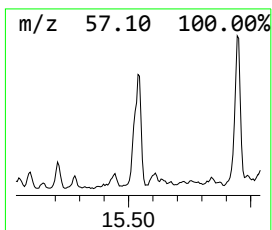
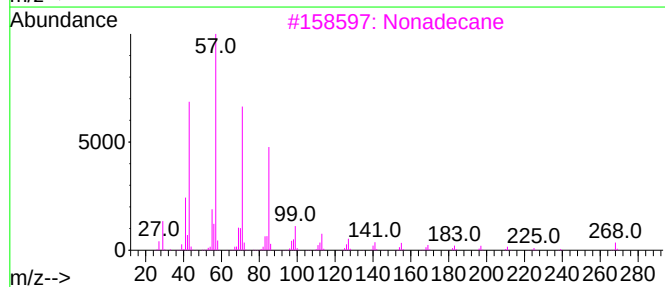
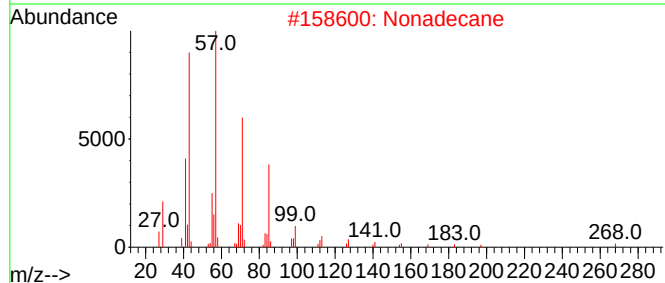
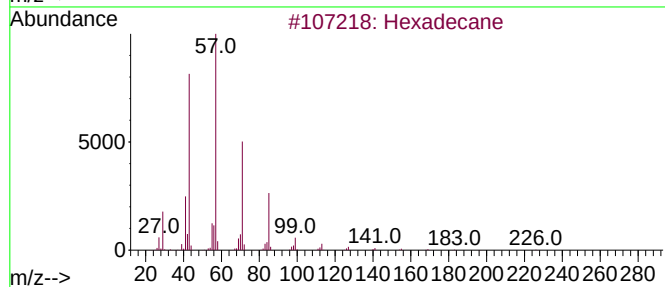
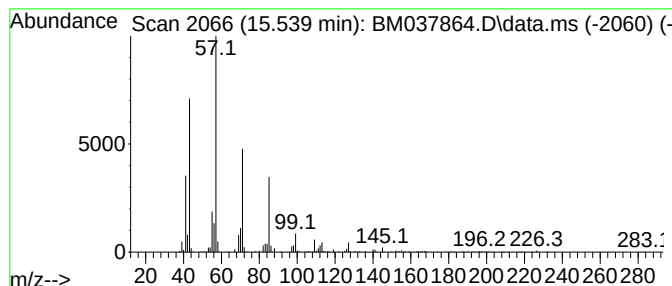
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
15.539	6.74 ng/ul	826663	Acenaphthene-d10		14.721	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexadecane		226	C16H34	000544-76-3	96
2	Nonadecane		268	C19H40	000629-92-5	80
3	Nonadecane		268	C19H40	000629-92-5	80
4	Dodecane, 1-iodo-		296	C12H25I	004292-19-7	72
5	Decane, 2,3,5-trimethyl-		184	C13H28	062238-11-3	72



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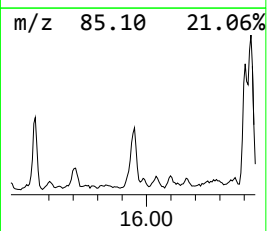
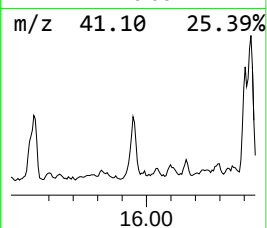
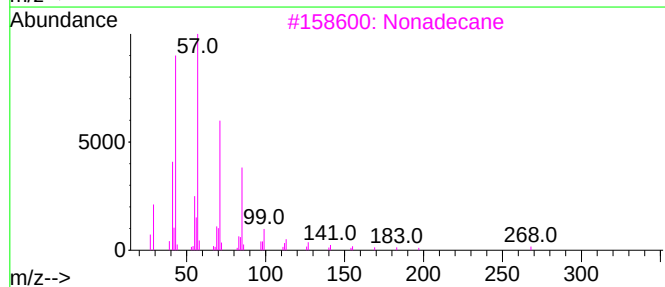
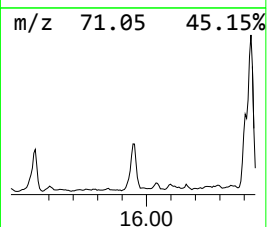
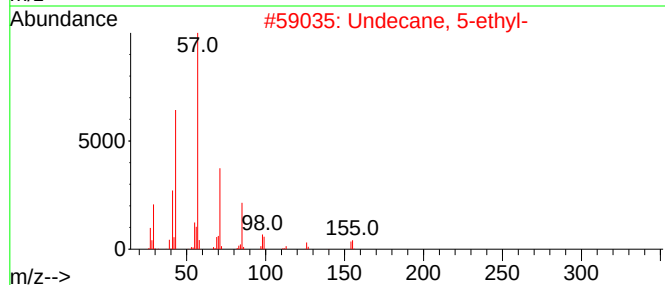
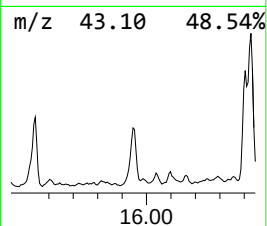
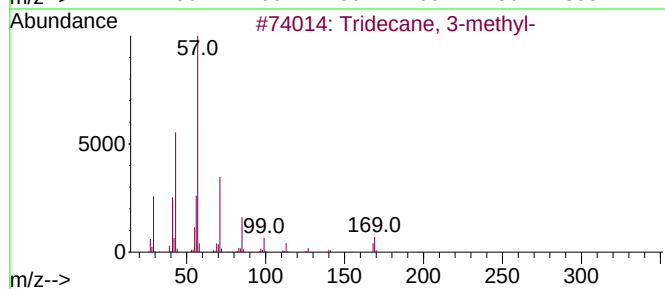
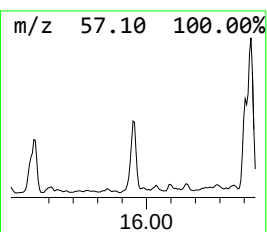
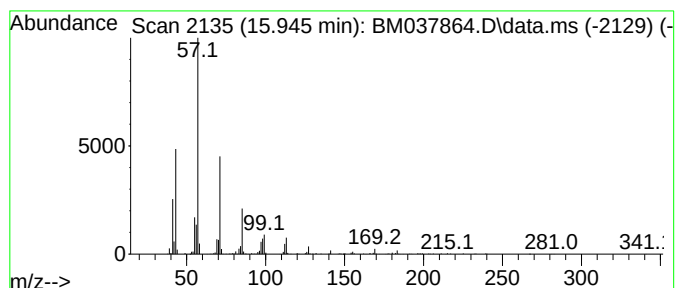
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Straight-Chai... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	5.80 ng/ul	710747	Acenaphthene-d10	14.721

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tridecane, 3-methyl-	198	C14H30	006418-41-3	87
2			Undecane, 5-ethyl-	184	C13H28	017453-94-0	78
3			Nonadecane	268	C19H40	000629-92-5	72
4			Heptadecane	240	C17H36	000629-78-7	64
5			Undecane, 3,5-dimethyl-	184	C13H28	017312-81-1	64



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GBNJ1MEDL

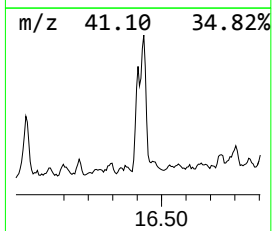
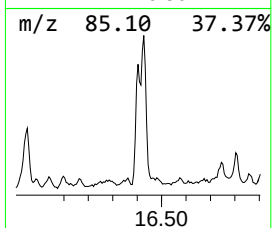
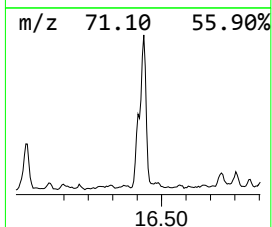
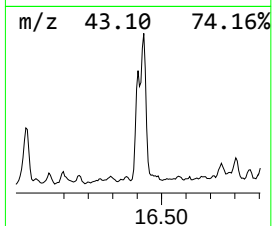
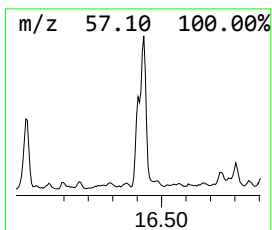
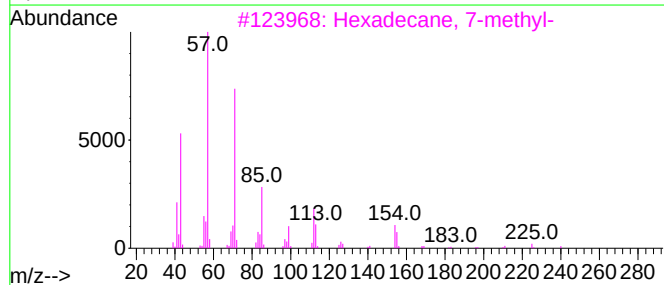
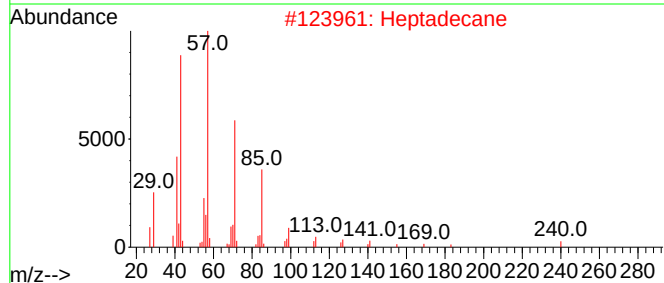
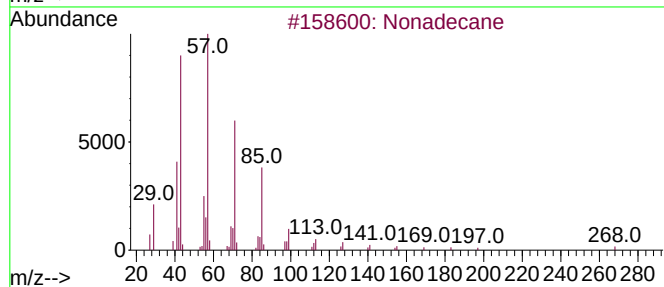
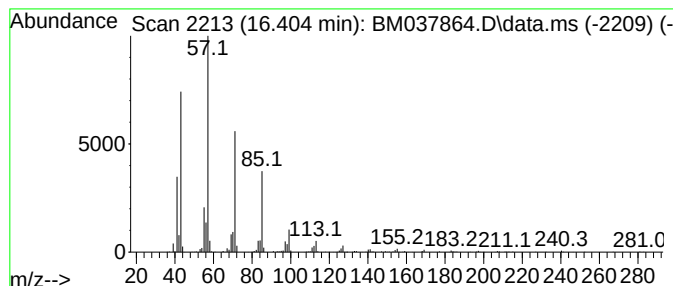
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.404	4.91 ng/ul	766055	Phenanthrene-d10	17.456

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonadecane	268	C19H40	000629-92-5	91
2			Heptadecane	240	C17H36	000629-78-7	91
3			Hexadecane, 7-methyl-	240	C17H36	026730-20-1	90
4			Tridecane	184	C13H28	000629-50-5	90
5			Tetracosane	338	C24H50	000646-31-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
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Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
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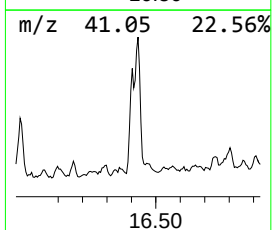
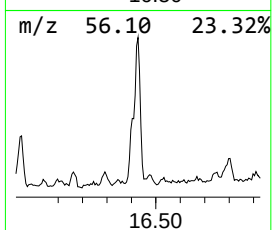
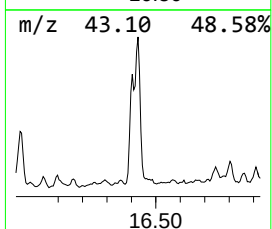
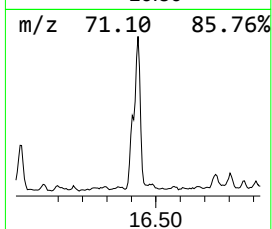
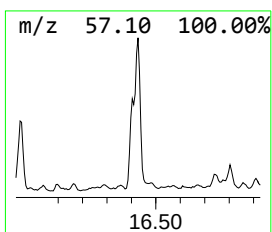
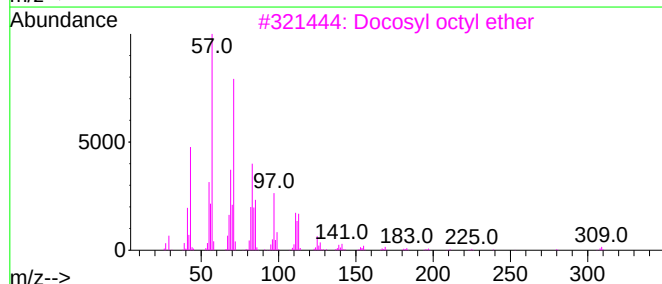
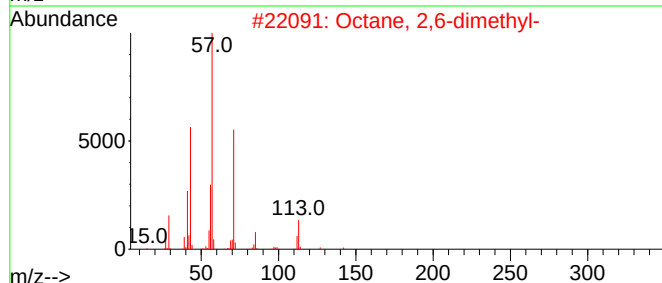
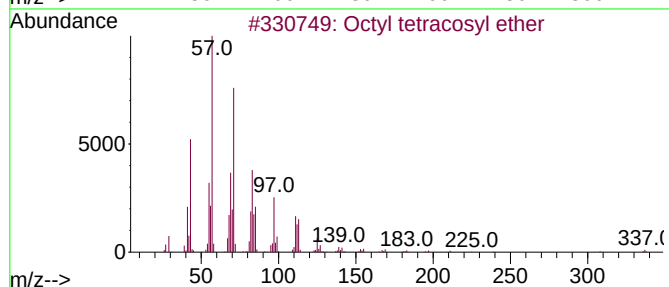
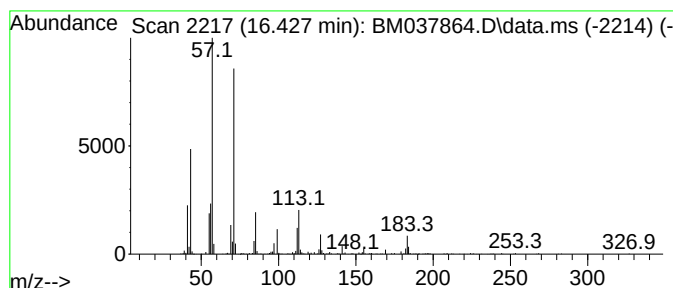
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Octyl tetraicosyl ether Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.427	9.71 ng/ul	1515030	Phenanthrene-d10	17.456

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octyl tetraicosyl ether	467	C32H66O	1000406-39-0	86
2			Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	81
3			Docosyl octyl ether	438	C30H62O	1000406-38-9	78
4			Sulfurous acid, 2-ethylhexyl non...	320	C17H36O3S	1010309-19-2	72
5			Sulfurous acid, 2-ethylhexyl und...	348	C19H40O3S	1000309-19-4	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
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Misc :
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Instrument :
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ClientSampleId :
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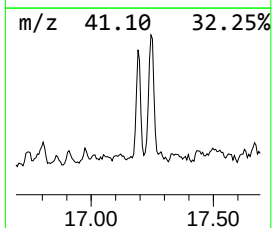
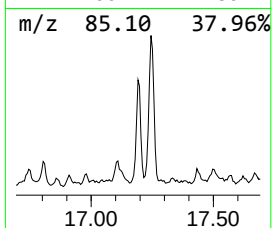
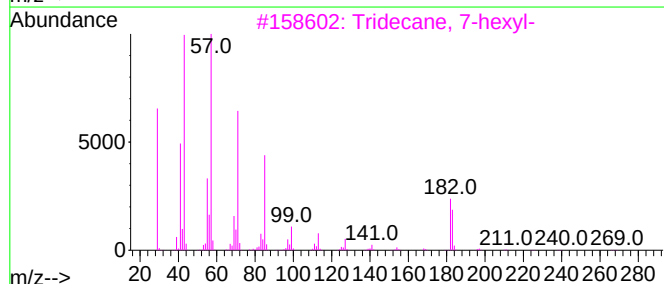
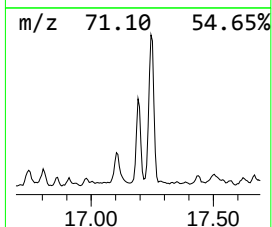
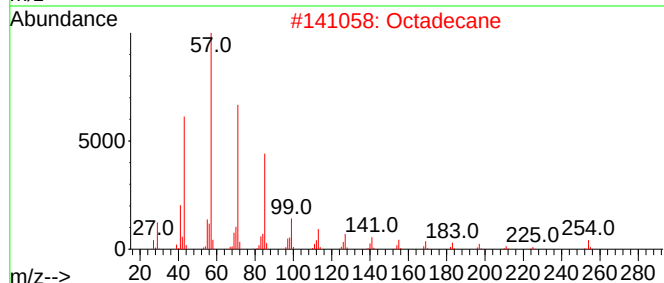
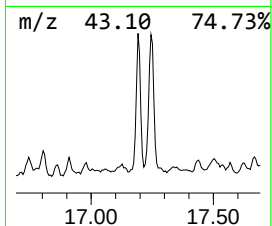
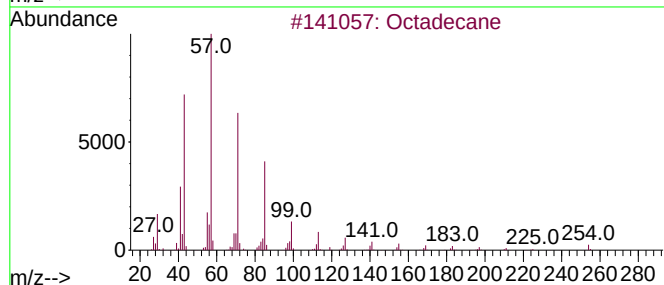
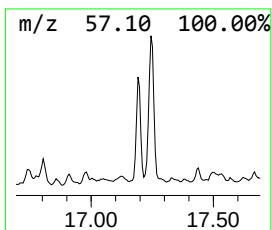
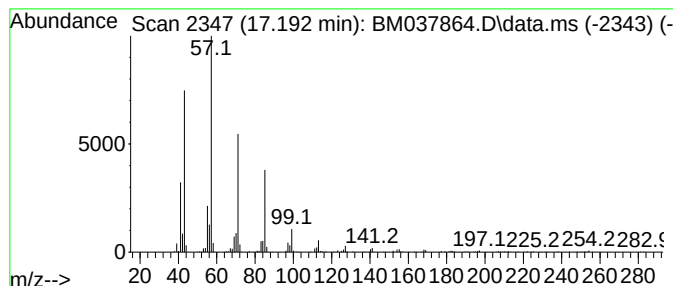
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.192	5.45 ng/ul	849877	Phenanthrene-d10	17.456

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Octadecane	254	C18H38	000593-45-3	93
2		Octadecane	254	C18H38	000593-45-3	90
3		Tridecane, 7-hexyl-	268	C19H40	007225-66-3	90
4		Hexadecane	226	C16H34	000544-76-3	90
5		Nonadecane	268	C19H40	000629-92-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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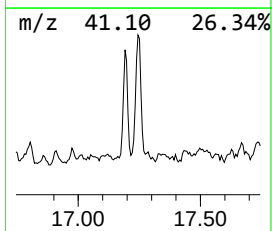
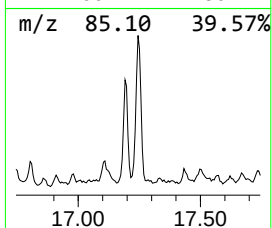
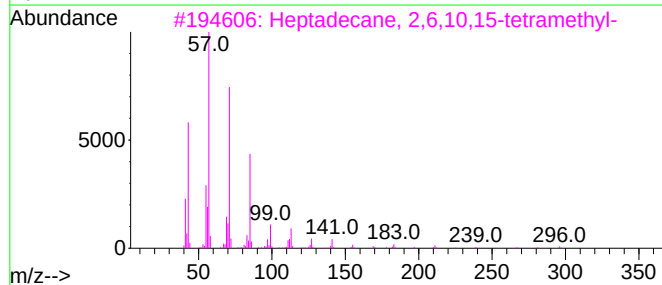
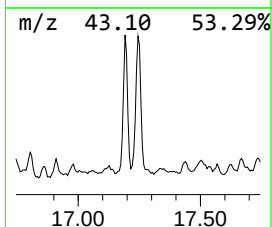
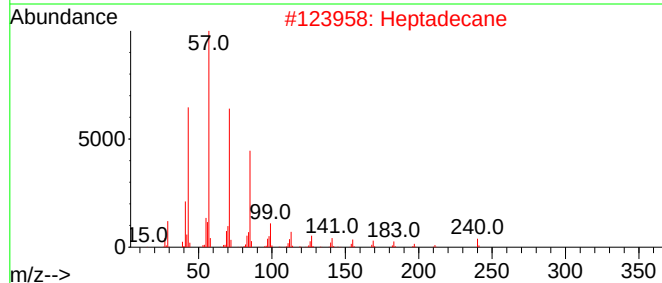
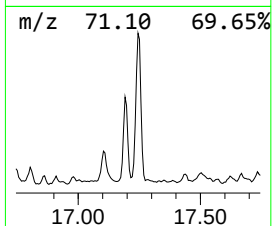
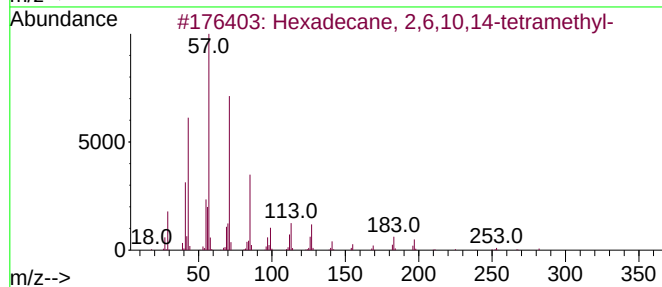
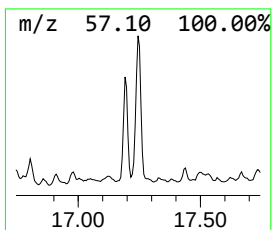
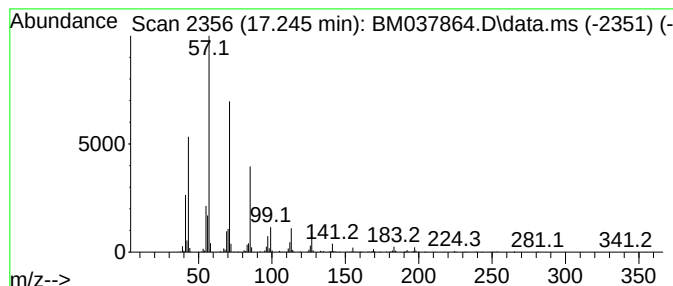
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
17.245	9.05 ng/ul	1412360	Phenanthrene-d10	17.456

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Hexadecane, 2,6,10,14-tetramethyl-	282	C20H42	000638-36-8	91
2		Heptadecane	240	C17H36	000629-78-7	90
3		Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	90
4		Tridecane, 2-methyl-	198	C14H30	001560-96-9	87
5		Tetracosane	338	C24H50	000646-31-1	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
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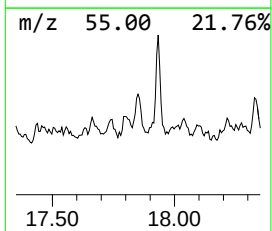
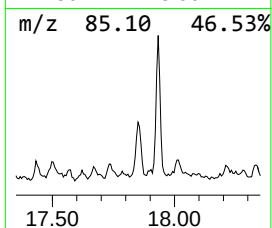
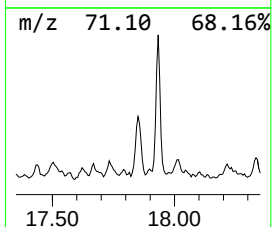
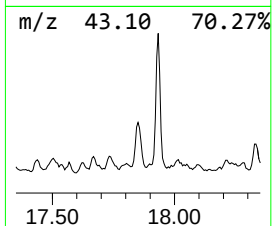
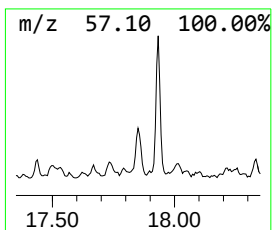
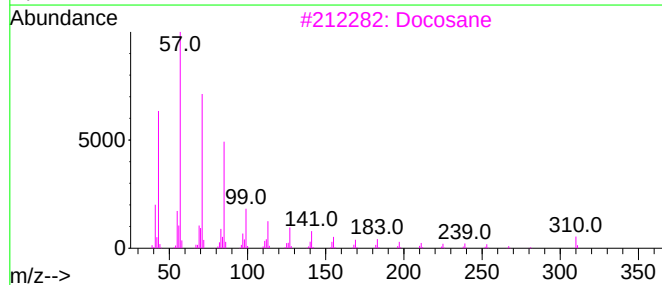
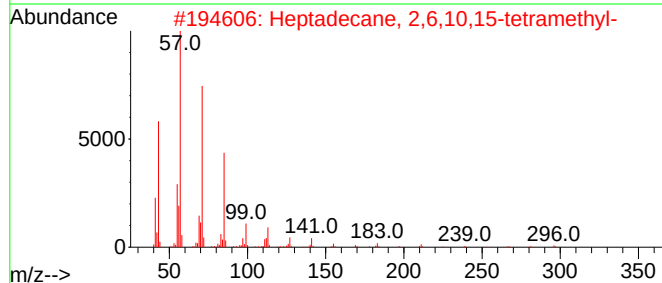
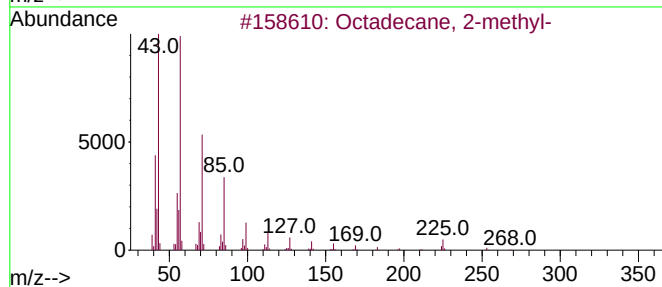
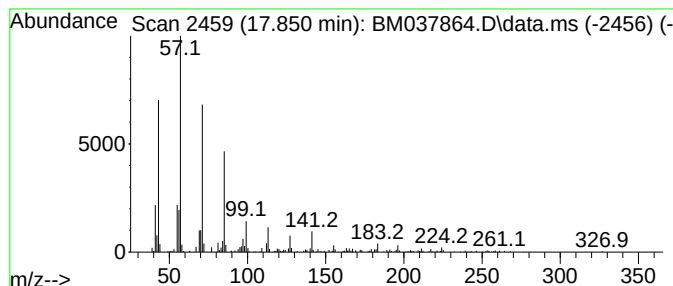
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.851	2.69 ng/ul	420180	Phenanthrene-d10	17.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Octadecane, 2-methyl-	268	C19H40	001560-88-9	90
2	Heptadecane, 2,6,10,15-tetramethyl-	296	C21H44	054833-48-6	87
3	Docosane	310	C22H46	000629-97-0	83
4	2-Bromotetradecane	276	C14H29Br	074036-95-6	80
5	Docosane, 1-iodo-	436	C22H45I	1000406-31-9	80



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
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Misc :
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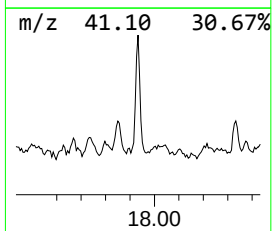
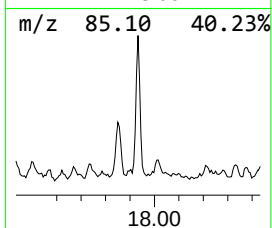
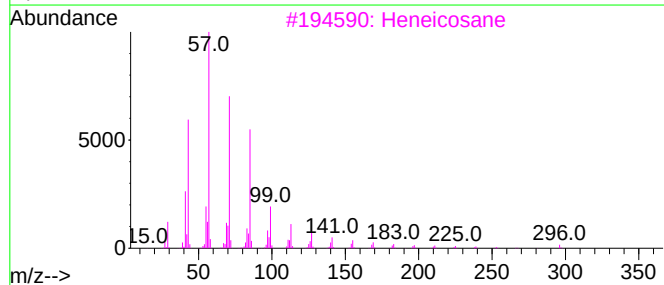
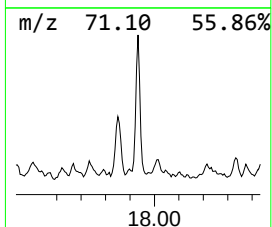
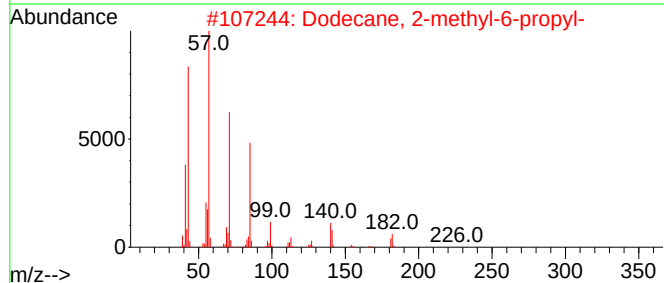
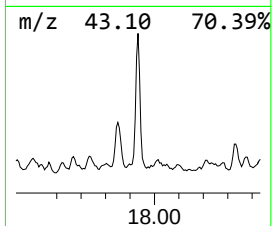
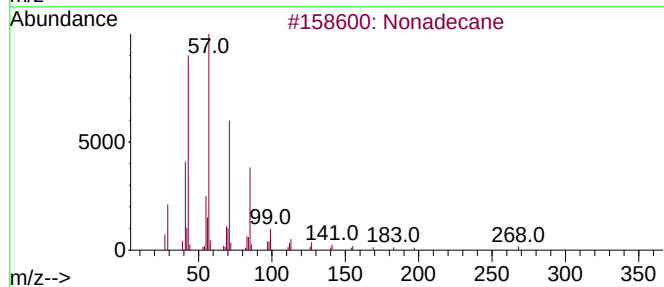
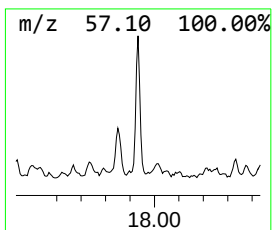
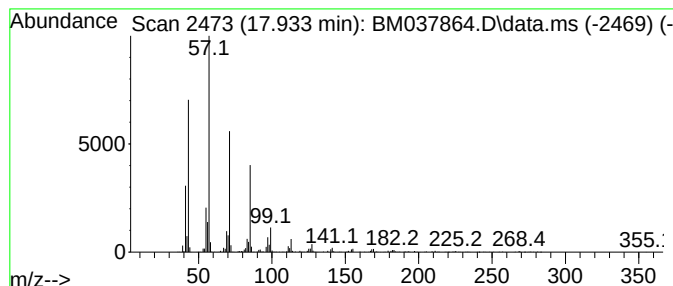
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
17.933	5.95 ng/ul	928322	Phenanthrene-d10		17.456	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Nonadecane		268	C19H40	000629-92-5	91
2	Dodecane, 2-methyl-6-propyl-		226	C16H34	055045-08-4	90
3	Heneicosane		296	C21H44	000629-94-7	90
4	Tetracosane		338	C24H50	000646-31-1	90
5	Eicosane		282	C20H42	000112-95-8	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
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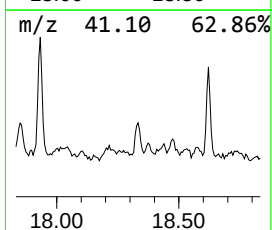
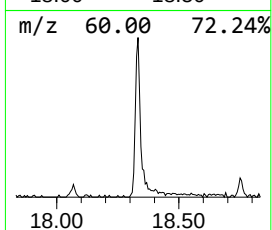
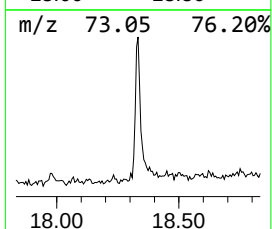
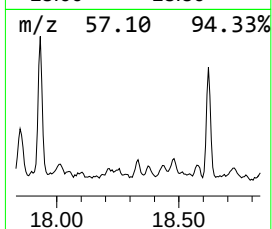
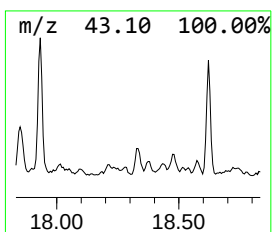
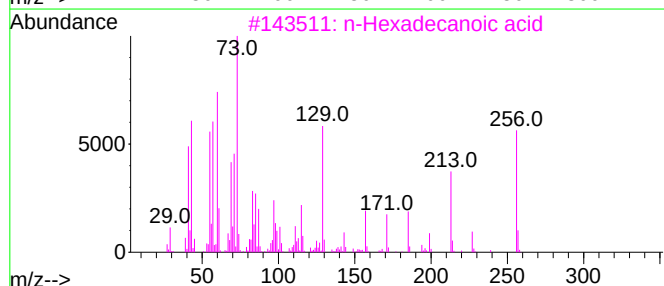
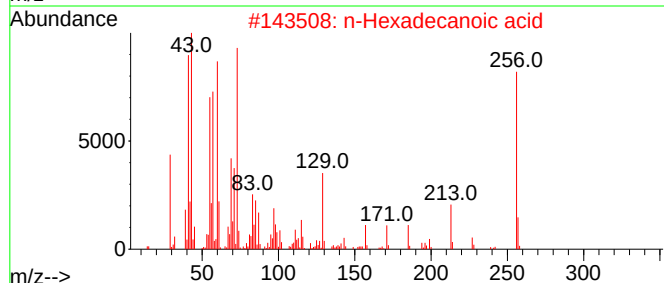
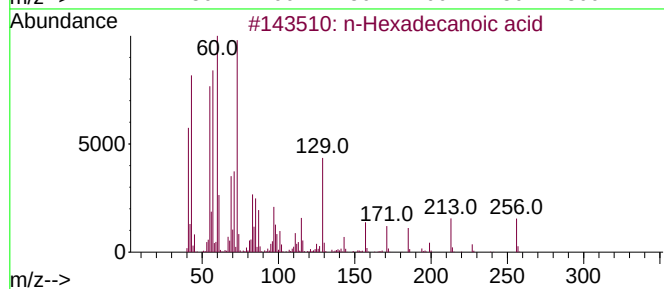
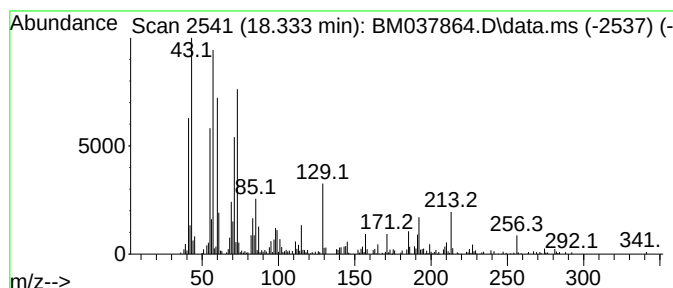
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.333	2.19 ng/ul	341625	Phenanthrene-d10	17.456	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
2	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	83
3	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	64
4	n-Hexadecanoic acid	256	C16H32O2	000057-10-3	60
5	Tetradecanoic acid	228	C14H28O2	000544-63-8	60



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Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ1MEDL

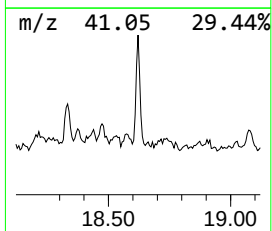
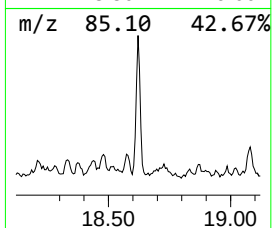
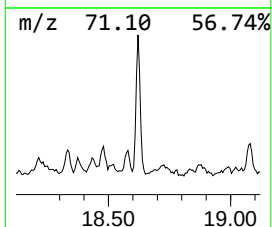
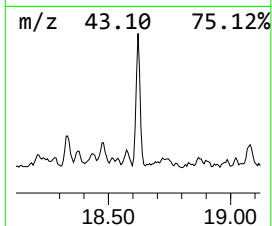
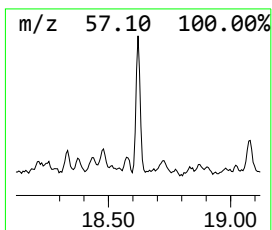
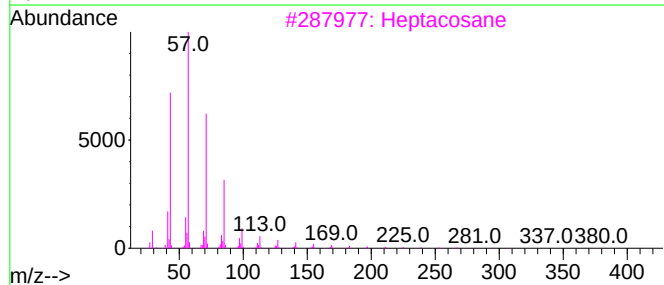
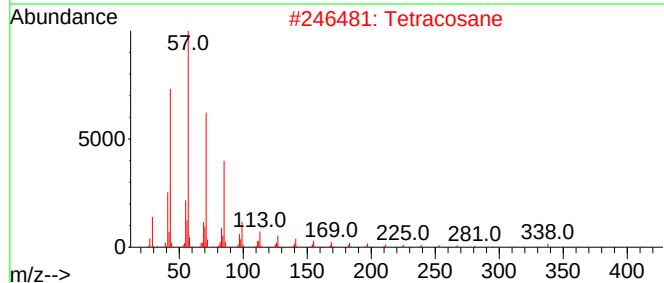
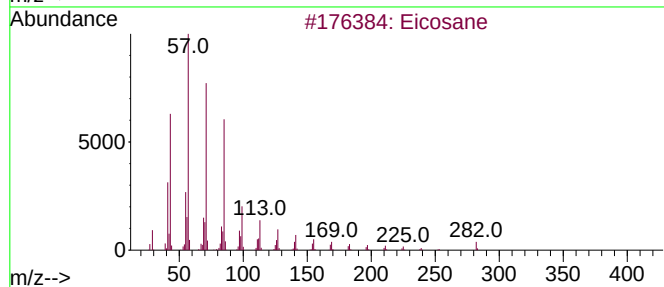
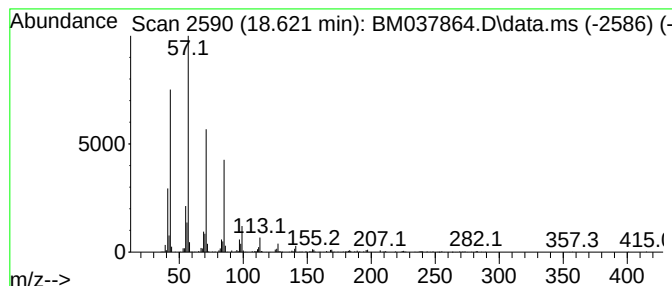
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Straight-Chai... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.621	4.50 ng/ul	701490	Phenanthrene-d10	17.456

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane	282	C20H42	000112-95-8	94
2		Tetracosane	338	C24H50	000646-31-1	91
3		Heptacosane	380	C27H56	000593-49-7	91
4		Pentadecane	212	C15H32	000629-62-9	91
5		Heneicosane	296	C21H44	000629-94-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

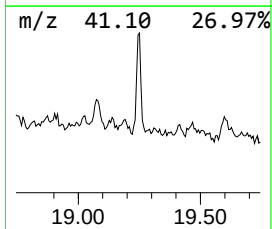
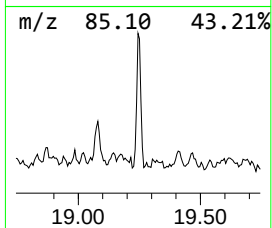
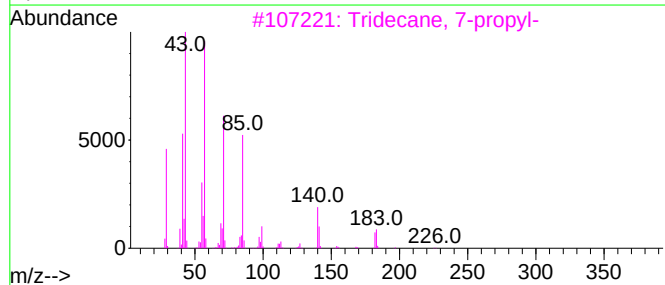
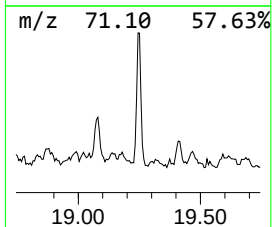
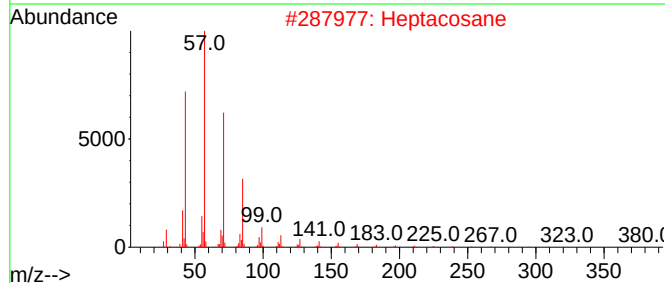
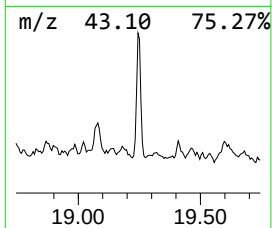
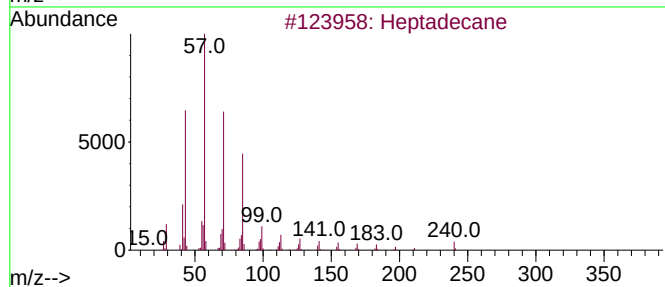
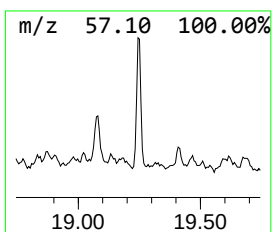
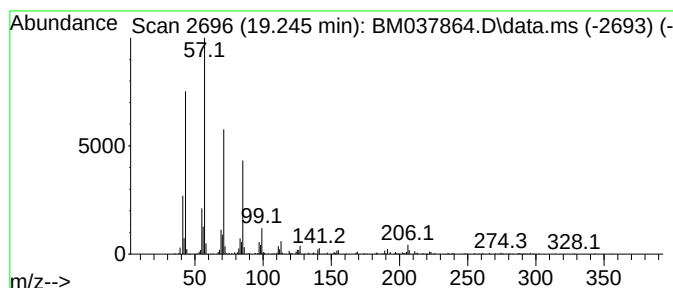
Instrument :
BNA_M
ClientSampleId :
GBNJ1MEDL

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 12 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.244	3.29 ng/ul	513590	Phenanthrene-d10		17.456	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane		240	C17H36	000629-78-7	93
2	Heptacosane		380	C27H56	000593-49-7	91
3	Tridecane, 7-propyl-		226	C16H34	055045-09-5	90
4	Nonadecane		268	C19H40	000629-92-5	90
5	Eicosane, 10-methyl-		296	C21H44	054833-23-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ1MEDL

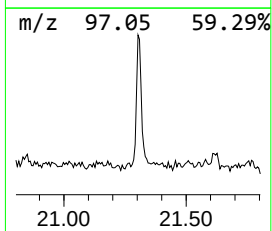
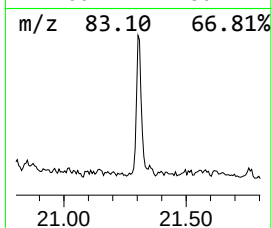
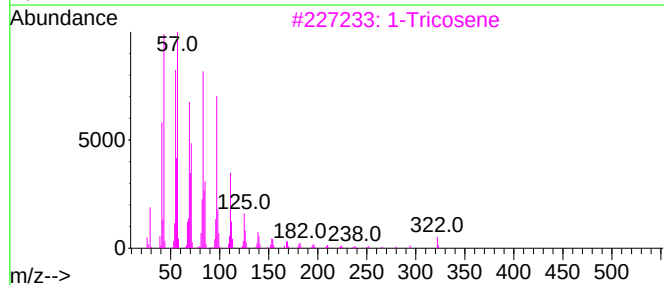
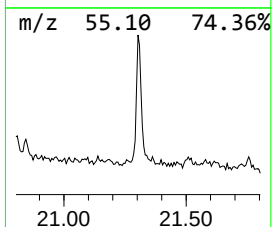
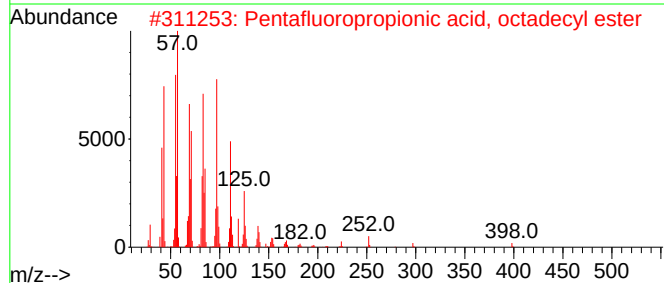
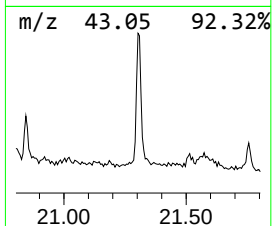
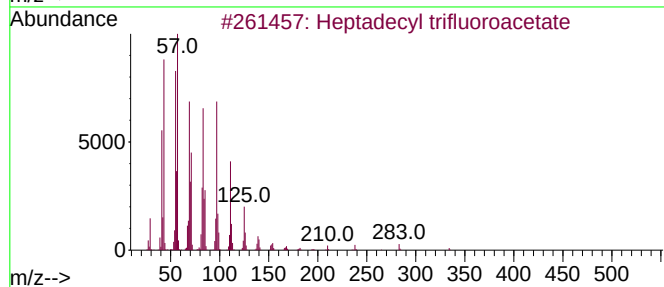
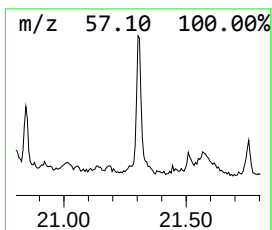
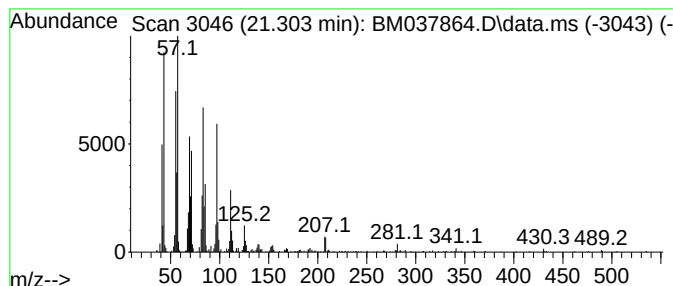
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM120122.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 Heptadecyl trifluoroacetate Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.303	3.55 ng/ul	529607	Chrysene-d12	21.621

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecyl trifluoroacetate	352	C19H35F3O2	1010351-87-0	93
2			Pentafluoropropionic acid, octadecyl ester	416	C21H37F5O2	959261-25-7	90
3			1-Tricosene	322	C23H46	018835-32-0	90
4			Pentafluoropropionic acid, pentafluorooctadecyl ester	374	C18H31F5O2	959092-08-1	90
5			Cyclooctacosane	392	C28H56	000297-24-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM120622\
Data File : BM037864.D
Acq On : 06 Dec 2022 21:12
Operator : CG/JU
Sample : N5738-12MEDL 2X
Misc :
ALS Vial : 15 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
GBNJ1MEDL

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

					--Internal Standard--			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
(DEL) Alkane: S...	14.598	9.6	ng/u1	1175910	3	14.721	2451760	20.0
(DEL) Alkane: S...	15.539	6.7	ng/u1	826663	3	14.721	2451760	20.0
(DEL) Alkane: S...	15.945	5.8	ng/u1	710747	3	14.721	2451760	20.0
(DEL) Alkane: S...	16.404	4.9	ng/u1	766055	4	17.456	3119760	20.0
Octyl tetracosy...	16.427	9.7	ng/u1	1515030	4	17.456	3119760	20.0
(DEL) Alkane: S...	17.192	5.5	ng/u1	849877	4	17.456	3119760	20.0
(DEL) Alkane: S...	17.245	9.1	ng/u1	1412360	4	17.456	3119760	20.0
(DEL) Alkane: S...	17.851	2.7	ng/u1	420180	4	17.456	3119760	20.0
(DEL) Alkane: S...	17.933	6.0	ng/u1	928322	4	17.456	3119760	20.0
n-Hexadecanoic ...	18.333	2.2	ng/u1	341625	4	17.456	3119760	20.0
(DEL) Alkane: S...	18.621	4.5	ng/u1	701490	4	17.456	3119760	20.0
(DEL) Alkane: S...	19.244	3.3	ng/u1	513590	4	17.456	3119760	20.0
Heptadecyl trif...	21.303	3.5	ng/u1	529607	5	21.621	2981230	20.0