

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
 Data File : BP015630.D
 Acq On : 20 Jun 2023 19:52
 Operator : MA/JU
 Sample : 03080-19
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFJ8

Integration Parameters: LSCINT.P

Integrator: RTE

Smoothing : OFF

Filtering: 5

Sampling : 1

Min Area: 0.1 % of largest Peak

Start Thrs: 0.2

Max Peaks: 100

Stop Thrs : 0

Peak Location: TOP

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

Peak separation: 5

Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015630.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.405	33	37	43	rVB	16682	22987	0.64%	0.057%
2	3.858	113	114	126	rVV	63245	105107	2.94%	0.261%
3	3.952	126	130	139	rVB	42253	62546	1.75%	0.155%
4	4.075	147	151	159	rVB2	13114	24336	0.68%	0.060%
5	4.175	165	168	176	rVB5	5601	10520	0.29%	0.026%
6	4.758	263	267	272	rVB6	6849	10697	0.30%	0.027%
7	5.240	346	349	357	rVB7	5154	9835	0.27%	0.024%
8	6.052	483	487	492	rVB6	6341	11625	0.32%	0.029%
9	6.299	524	529	539	rBV5	9404	24839	0.69%	0.062%
10	6.734	597	603	609	rVB4	19835	35679	1.00%	0.089%
11	7.046	649	656	661	rBV9	10429	20373	0.57%	0.051%
12	7.246	679	690	702	rBV	376970	656332	18.34%	1.630%
13	7.405	711	717	729	rVB	305795	490766	13.72%	1.219%
14	7.499	729	733	742	rVB3	14058	27515	0.77%	0.068%
15	7.611	744	752	761	rBV	416398	689012	19.26%	1.711%
16	7.675	761	763	768	rVV5	9874	14915	0.42%	0.037%
17	8.081	820	832	843	rBV	358500	603859	16.88%	1.500%
18	8.281	859	866	874	rBV10	7331	18610	0.52%	0.046%
19	8.352	874	878	882	rBV6	5534	10912	0.30%	0.027%
20	8.805	948	955	967	rBV	393139	692766	19.36%	1.721%
21	8.928	972	976	981	rVB6	5693	9609	0.27%	0.024%
22	9.263	1024	1033	1046	rBV	422020	744856	20.82%	1.850%
23	9.416	1056	1059	1067	rVB10	5603	11224	0.31%	0.028%
24	9.993	1147	1157	1165	rBV	361973	647167	18.09%	1.607%
25	10.534	1242	1249	1269	rBV	441950	904116	25.27%	2.246%
26	10.681	1269	1274	1283	rVB	5326	17328	0.48%	0.043%
27	10.857	1299	1304	1306	rBV3	9678	12709	0.36%	0.032%
28	10.910	1306	1313	1319	rVV	550118	895149	25.02%	2.223%
29	10.957	1319	1321	1326	rVB	15860	22461	0.63%	0.056%
30	11.057	1331	1338	1358	rBV	228096	527980	14.76%	1.311%
31	11.716	1445	1450	1454	rBV8	6360	12868	0.36%	0.032%
32	11.910	1477	1483	1487	rBV2	13707	22949	0.64%	0.057%
33	12.257	1535	1542	1545	rBV7	9632	16342	0.46%	0.041%
34	12.299	1546	1549	1555	rVB3	11143	16711	0.47%	0.042%
35	12.493	1577	1582	1588	rBV3	10045	18713	0.52%	0.046%
36	12.563	1590	1594	1601	rVB	14492	25181	0.70%	0.063%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	12.722	1611	1621	1626	rBV7	9686	26471	0.74%	0.066%
38	12.781	1626	1631	1639	rVB2	14282	27783	0.78%	0.069%
39	13.240	1705	1709	1713	rBV3	10900	16573	0.46%	0.041%
40	13.328	1716	1724	1731	rVB3	7768	19285	0.54%	0.048%
41	13.534	1752	1759	1762	rVV2	17630	26683	0.75%	0.066%
42	13.610	1766	1772	1781	rVV3	24135	48489	1.36%	0.120%
43	13.916	1822	1824	1830	rVB6	7406	9205	0.26%	0.023%
44	14.045	1841	1846	1851	rBV4	24264	39669	1.11%	0.099%
45	14.134	1852	1861	1867	rBV	1043029	1439494	40.23%	3.575%
46	14.187	1867	1870	1876	rVV3	20173	30159	0.84%	0.075%
47	14.251	1877	1881	1884	rVV3	14254	20829	0.58%	0.052%
48	14.422	1903	1910	1924	rBV	1151176	1769773	49.46%	4.396%
49	14.545	1927	1931	1935	rVB7	5929	9078	0.25%	0.023%
50	14.598	1935	1940	1947	rVB	24149	32391	0.91%	0.080%
51	14.675	1949	1953	1956	rBV3	16108	27915	0.78%	0.069%
52	14.728	1956	1962	1969	rVB	852890	1231401	34.41%	3.058%
53	14.957	1995	2001	2021	rBV	147748	446262	12.47%	1.108%
54	15.128	2023	2030	2035	rVV7	29606	61957	1.73%	0.154%
55	15.340	2062	2066	2069	rBV6	8783	13469	0.38%	0.033%
56	15.398	2073	2076	2080	rVB5	7838	12087	0.34%	0.030%
57	15.504	2090	2094	2098	rBV7	13637	21935	0.61%	0.054%
58	15.551	2098	2102	2106	rVB2	27999	38300	1.07%	0.095%
59	15.722	2124	2131	2137	rBV	1474033	2128031	59.47%	5.285%
60	15.769	2137	2139	2148	rVV4	31701	49983	1.40%	0.124%
61	15.857	2148	2154	2166	rVV	359770	557061	15.57%	1.384%
62	15.957	2167	2171	2178	rVB7	18286	32249	0.90%	0.080%
63	16.087	2187	2193	2204	rVB	290454	432419	12.09%	1.074%
64	16.434	2245	2252	2261	rBV2	47089	113575	3.17%	0.282%
65	16.575	2271	2276	2282	rBV7	15136	25348	0.71%	0.063%
66	16.857	2321	2324	2331	rVB3	17820	26580	0.74%	0.066%
67	17.016	2346	2351	2355	rBV7	19869	35296	0.99%	0.088%
68	17.139	2369	2372	2377	rBV2	16949	29390	0.82%	0.073%
69	17.210	2380	2384	2390	rBV4	46382	77609	2.17%	0.193%
70	17.357	2405	2409	2416	rVB	66292	91850	2.57%	0.228%
71	17.422	2417	2420	2425	rBV2	49691	60244	1.68%	0.150%
72	17.486	2425	2431	2435	rBV	1145235	1627758	45.49%	4.043%
73	17.528	2435	2438	2442	rVV	106597	144316	4.03%	0.358%
74	17.586	2442	2448	2459	rVB	1633061	2299701	64.27%	5.712%
75	17.951	2506	2510	2514	rBV	39209	49047	1.37%	0.122%
76	18.175	2544	2548	2553	rBV3	57613	90196	2.52%	0.224%

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77	18.228	2553	2557	2562	rBV2	41528	60149	1.68%	0.149%
78	18.286	2563	2567	2572	rBV2	80784	110264	3.08%	0.274%
79	18.339	2572	2576	2582	rBV	847969	1191278	33.29%	2.959%
80	18.528	2604	2608	2612	rBV3	41244	63020	1.76%	0.157%
81	18.616	2620	2623	2631	rBV2	59211	102562	2.87%	0.255%
82	19.222	2723	2726	2730	rVB2	90880	99125	2.77%	0.246%
83	19.304	2738	2740	2746	rVB5	56925	98878	2.76%	0.246%
84	19.428	2758	2761	2763	rBV3	62656	82929	2.32%	0.206%
85	19.480	2766	2770	2777	rVV	402497	571342	15.97%	1.419%
86	19.569	2781	2785	2790	rVV	289431	426466	11.92%	1.059%
87	19.616	2790	2793	2799	rVB6	109693	156190	4.37%	0.388%
88	19.775	2816	2820	2821	rBV3	158111	186032	5.20%	0.462%
89	19.810	2821	2826	2838	rVB2	2336647	3578122	100.00%	8.887%
90	20.292	2906	2908	2914	rVB3	113362	158587	4.43%	0.394%
91	20.680	2972	2974	2978	rBV4	81995	96971	2.71%	0.241%
92	20.780	2988	2991	2995	rBV	97905	109913	3.07%	0.273%
93	21.239	3065	3069	3075	rVB2	604318	880977	24.62%	2.188%
94	21.569	3119	3125	3129	rBV	1453395	2136422	59.71%	5.306%
95	22.163	3223	3226	3230	rVB2	291741	342611	9.58%	0.851%
96	23.280	3411	3416	3421	rVB	691784	1086190	30.36%	2.698%
97	23.951	3524	3530	3536	rBV2	1476405	2862096	79.99%	7.109%
98	24.116	3552	3558	3565	rVB	863039	1702581	47.58%	4.229%
99	24.727	3656	3662	3675	rVB	958361	2133122	59.62%	5.298%
100	26.710	3994	3999	4012	rVB	513722	1337693	37.39%	3.322%

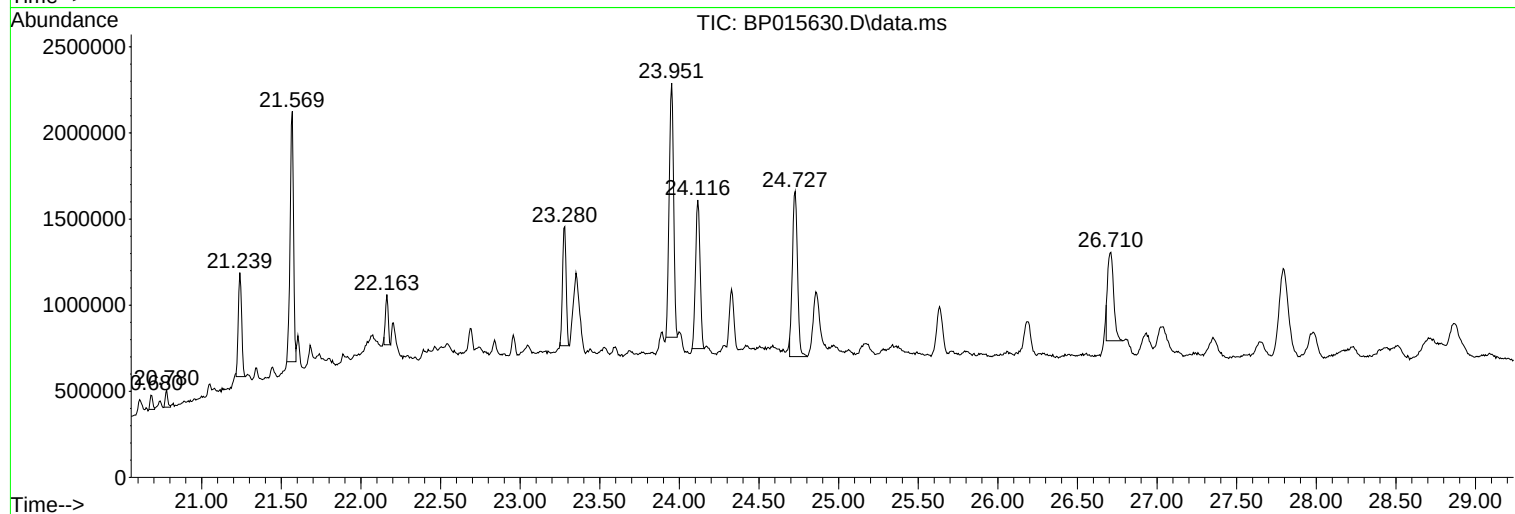
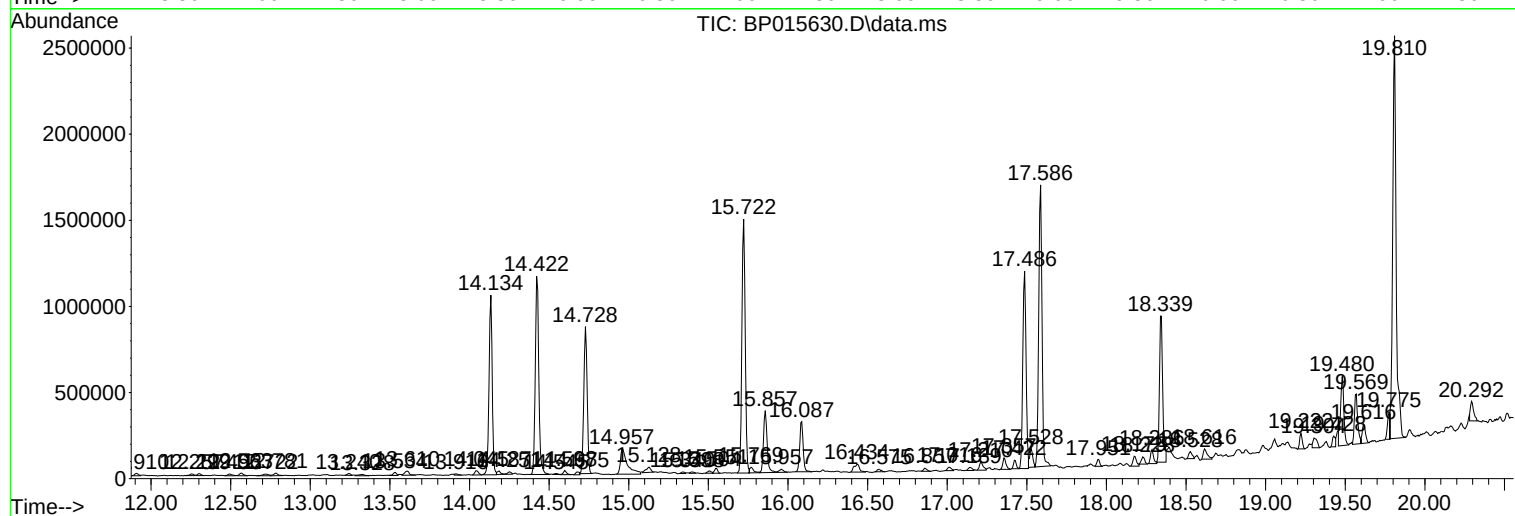
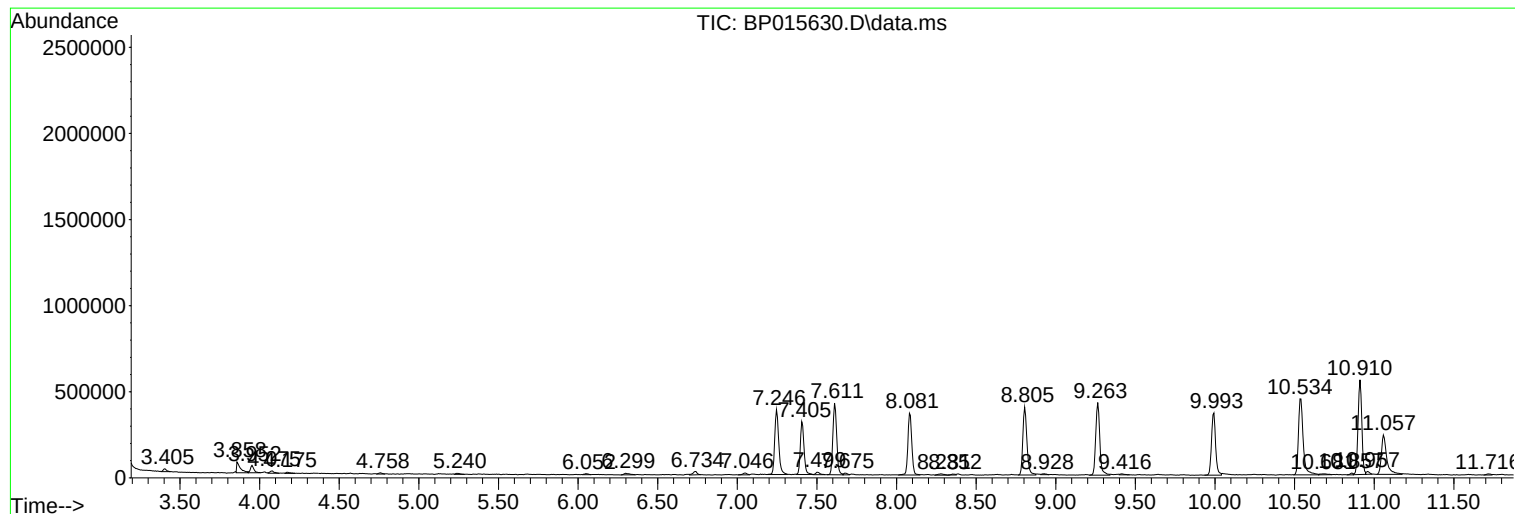
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Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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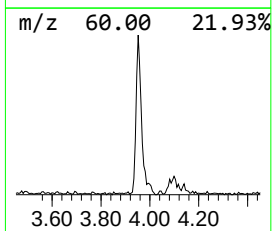
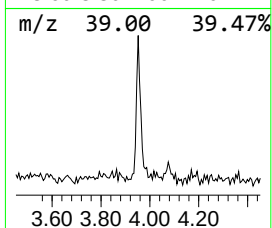
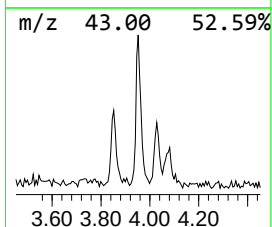
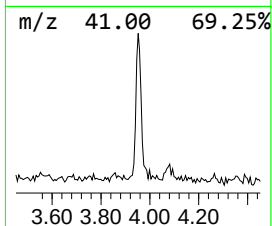
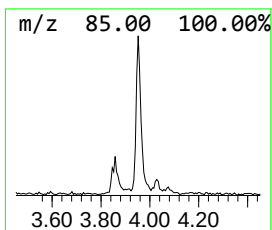
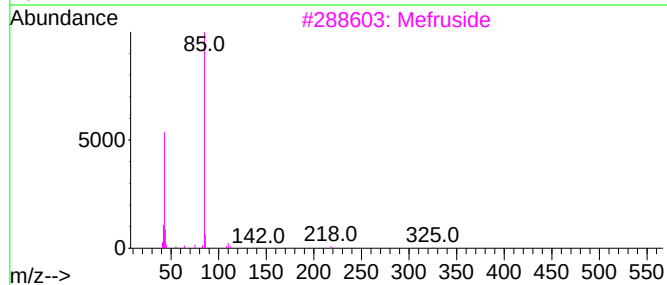
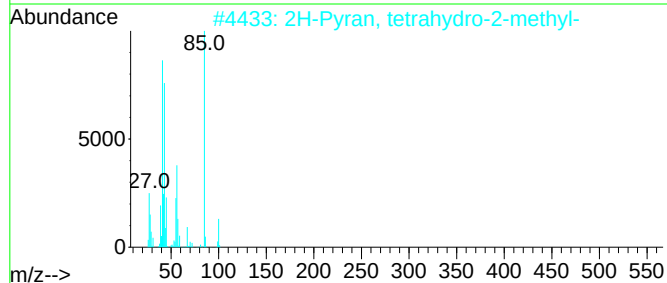
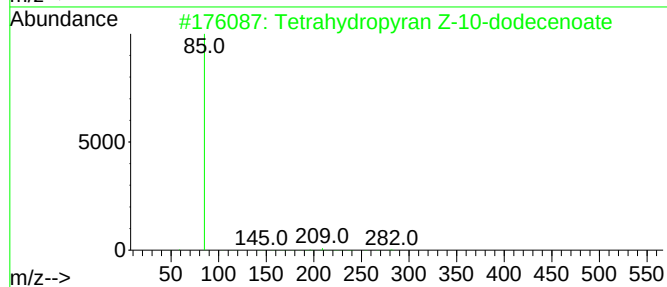
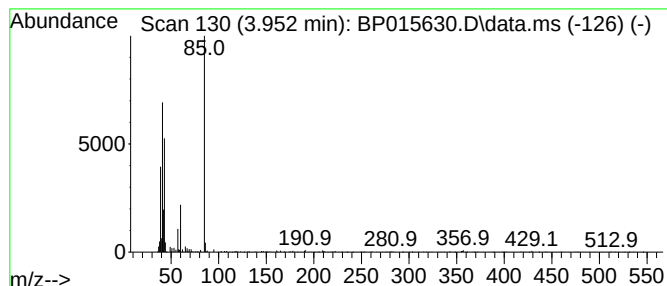
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrahydropyran Z-10-dodece... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.952	2.07 ng/ul	62546	1,4-Dichlorobenzene-d4	8.081

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrahydropyran Z-10-dodecenoate	282	C17H30O3	1000130-96-5	72
2			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	38
3			Mefruside	382	C13H19ClN2O5S2	007195-27-9	28
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	28
5			Oxirane, (2-methylpropyl)-	100	C6H12O	023850-78-4	25



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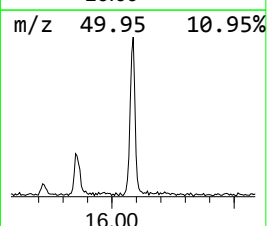
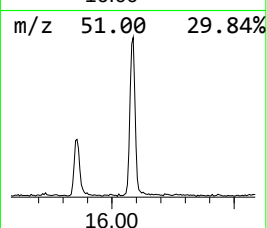
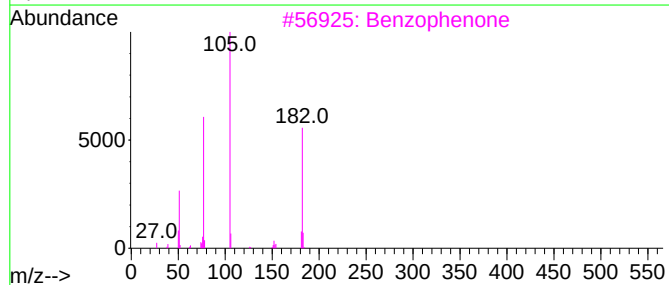
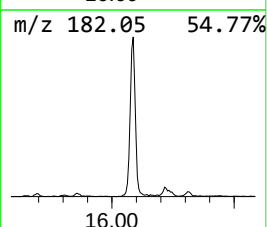
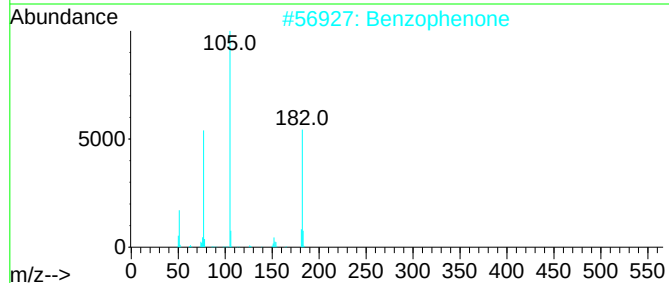
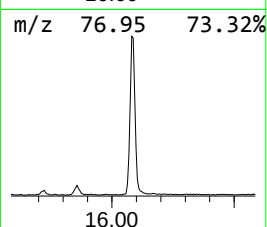
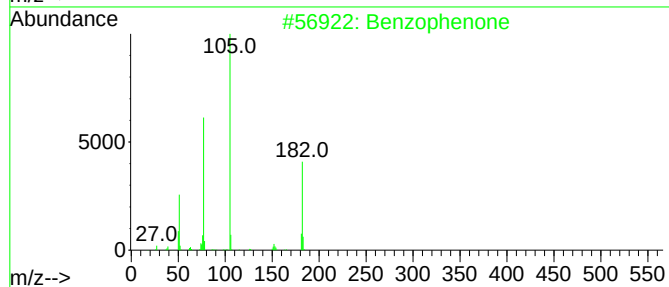
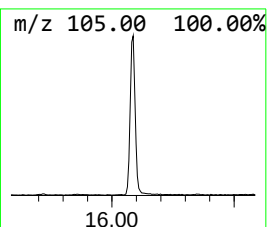
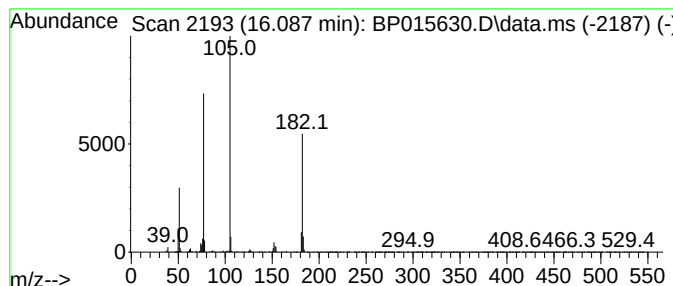
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.087	7.02 ng/ul	432419	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	97
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzophenone	182	C13H10O	000119-61-9	91



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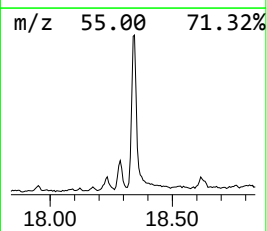
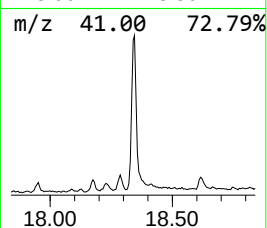
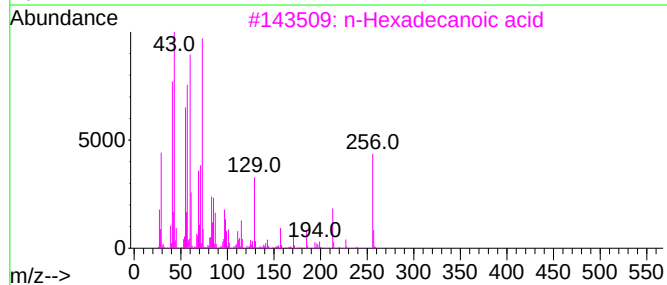
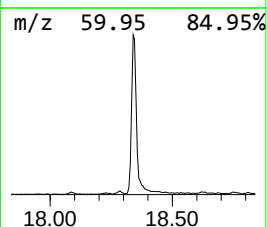
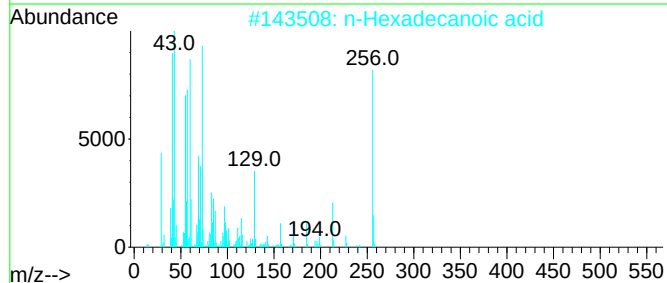
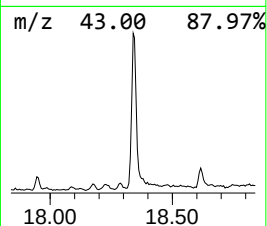
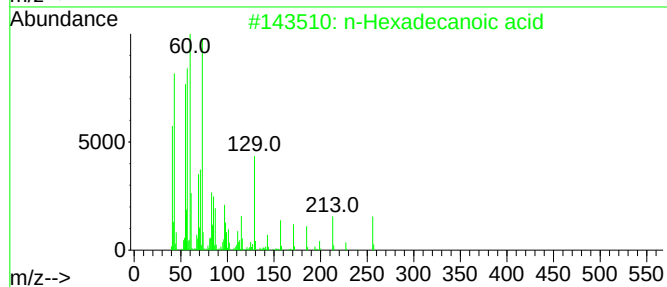
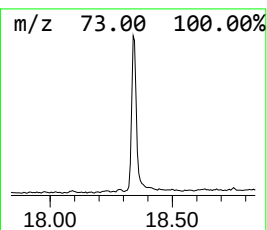
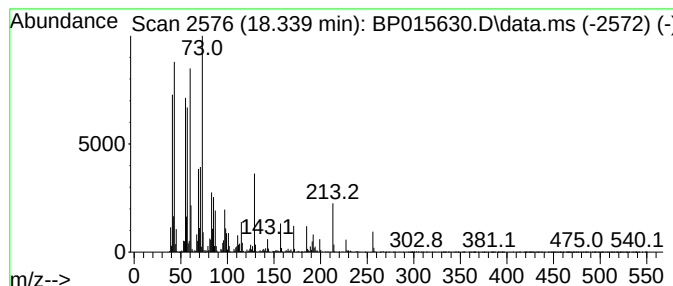
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.339	14.64 ng/ul	1191280	Phenanthrene-d10	17.486

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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
Acq On : 20 Jun 2023 19:52
Operator : MA/JU
Sample : 03080-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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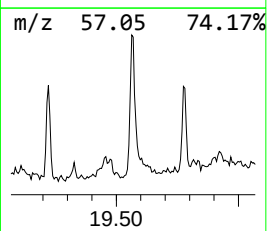
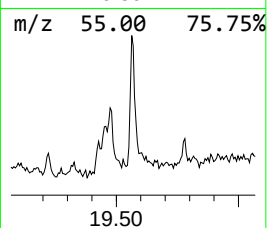
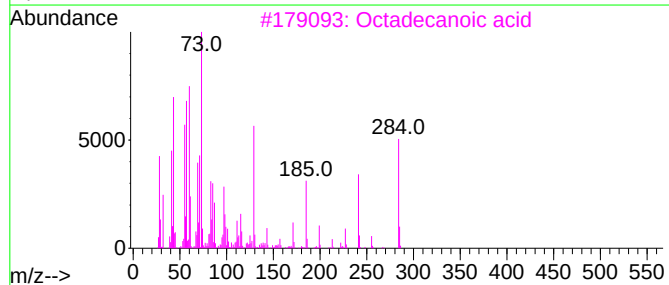
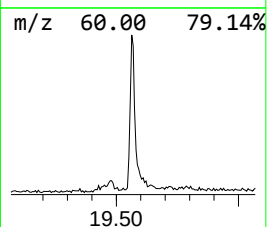
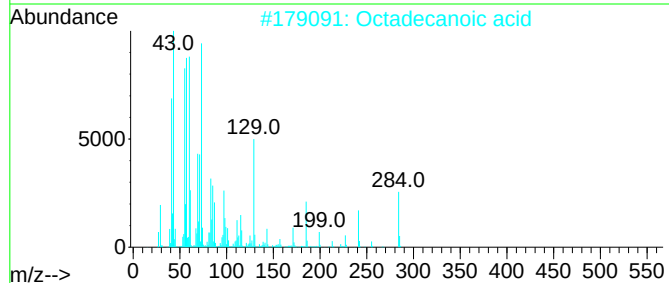
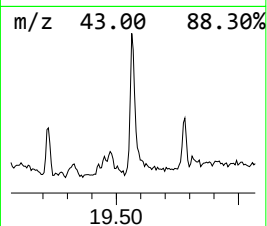
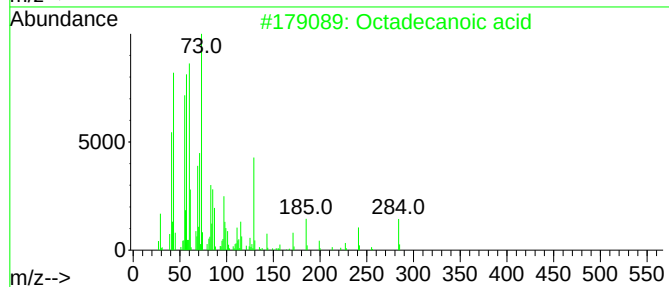
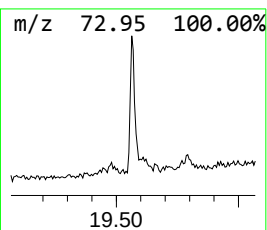
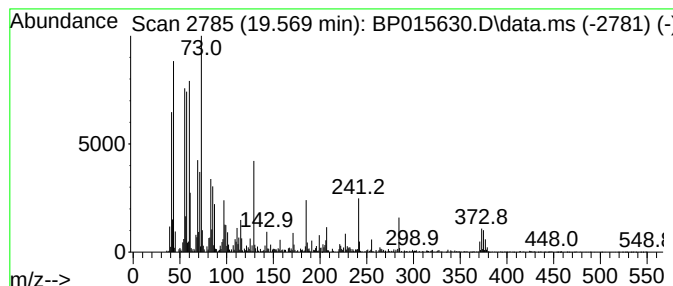
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.569	3.99 ng/ul	426466	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	95
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Pentadecanoic acid	242	C15H30O2	001002-84-2	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
Acq On : 20 Jun 2023 19:52
Operator : MA/JU
Sample : 03080-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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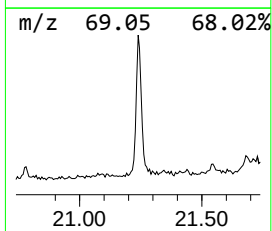
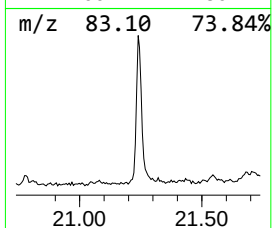
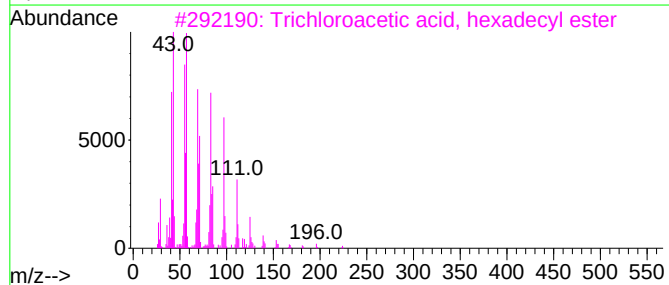
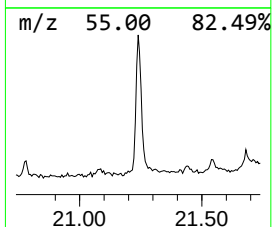
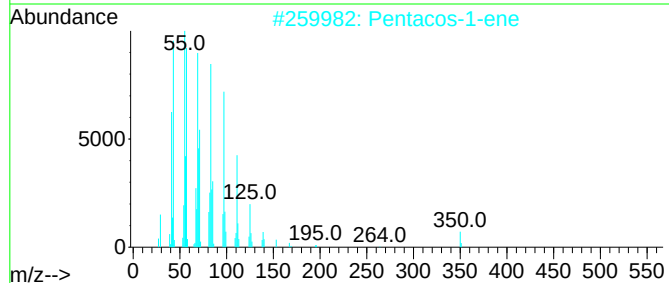
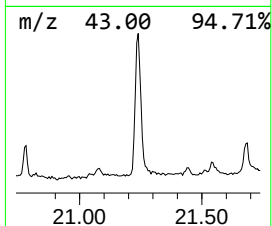
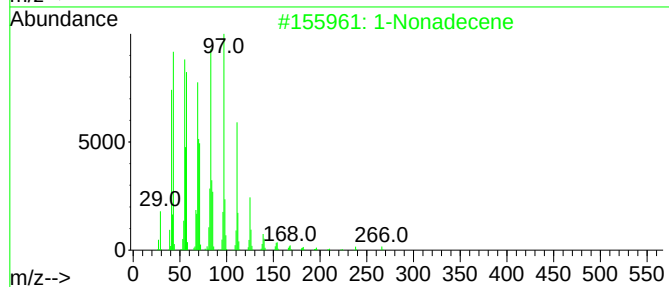
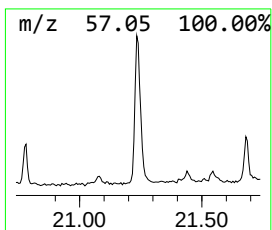
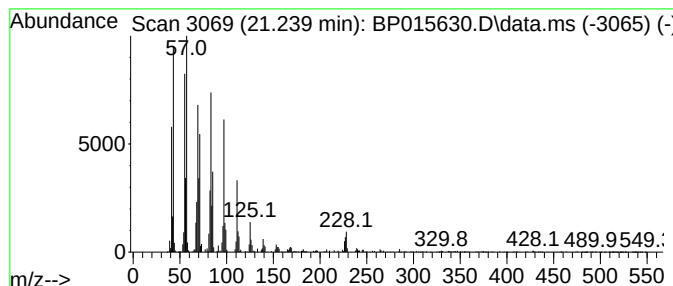
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.239	8.25 ng/ul	880977	Chrysene-d12	21.569

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Nonadecene	266	C19H38	018435-45-5	96
2			Pentacos-1-ene	350	C25H50	016980-85-1	95
3			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	93
4			1-Hexacosene	364	C26H52	018835-33-1	93
5			17-Pentatriacontene	491	C35H70	006971-40-0	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
Acq On : 20 Jun 2023 19:52
Operator : MA/JU
Sample : 03080-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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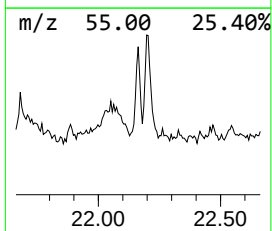
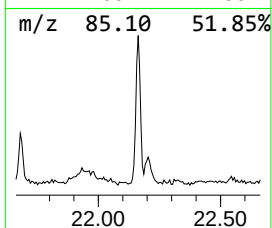
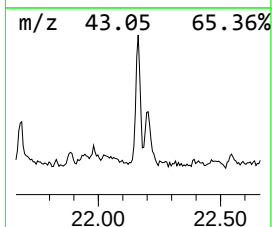
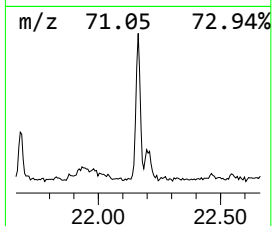
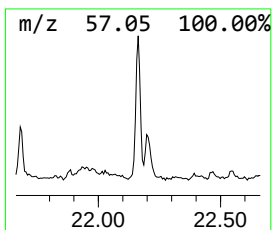
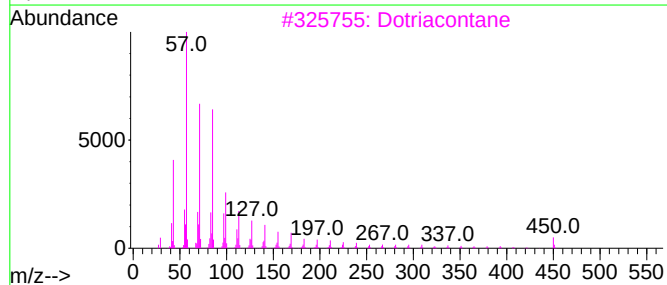
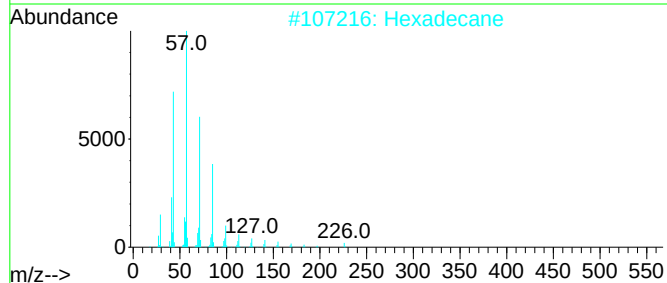
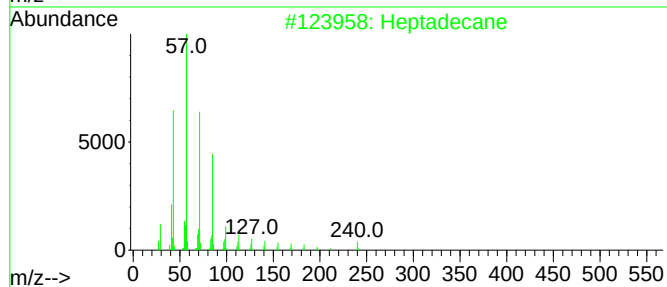
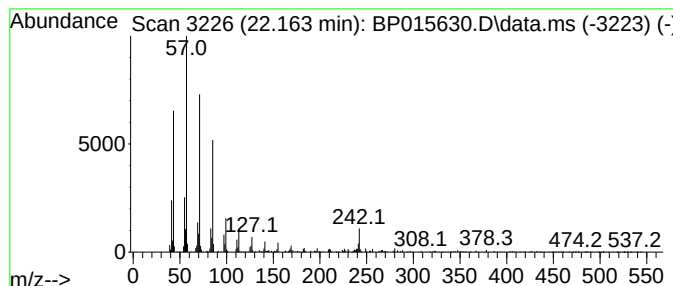
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.163	3.21 ng/ul	342611	Chrysene-d12	21.569

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Heptadecane	240	C17H36	000629-78-7	91
2		Hexadecane	226	C16H34	000544-76-3	91
3		Dotriacontane	451	C32H66	000544-85-4	87
4		Heptadecane	240	C17H36	000629-78-7	87
5		Hentriacontane	437	C31H64	000630-04-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
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Sample : 03080-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

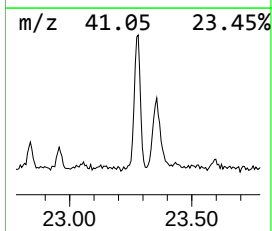
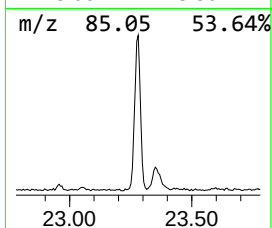
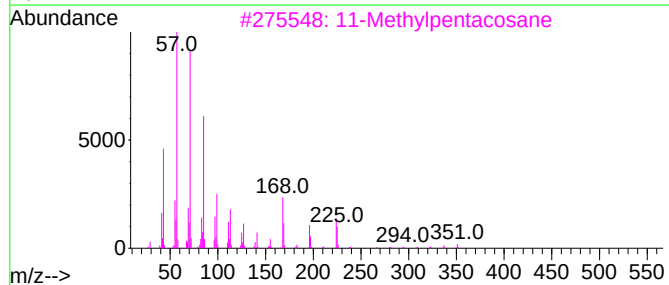
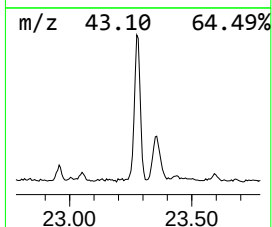
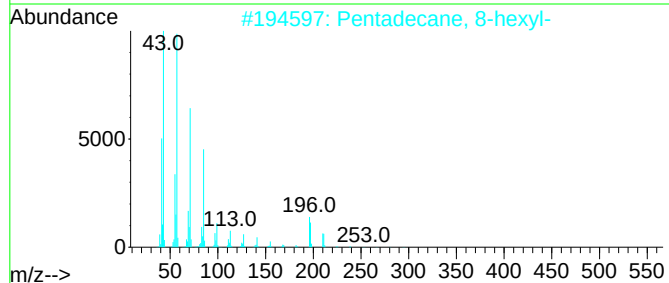
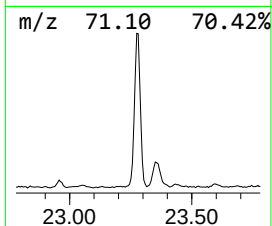
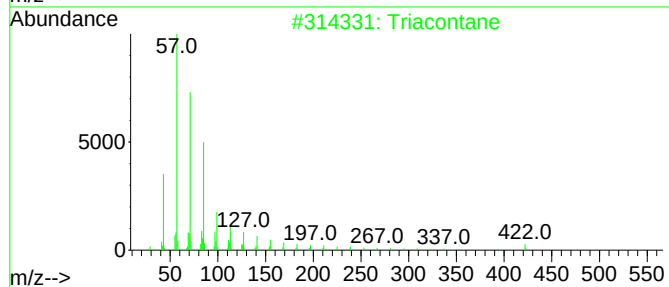
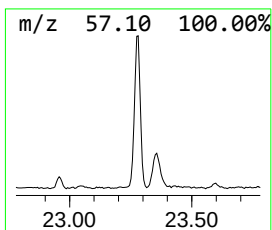
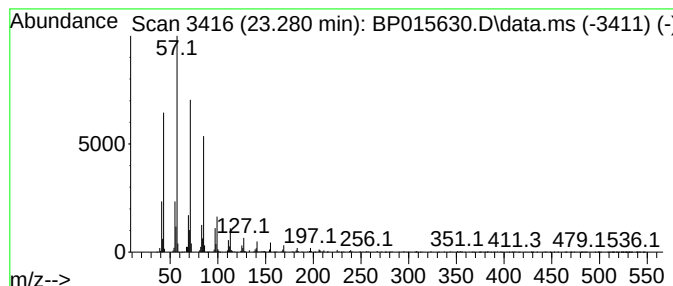
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD		R.T.	
23.280	12.76 ng/ul	1086190	Perylene-d12		24.116	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Triacontane		422	C30H62	000638-68-6	97
2	Pentadecane, 8-hexyl-		296	C21H44	013475-75-7	93
3	11-Methylpentacosane		366	C26H54	015689-71-1	93
4	Tetracosane		338	C24H50	000646-31-1	91
5	Octacosane, 2-methyl-		408	C29H60	001560-98-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
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Sample : 03080-19
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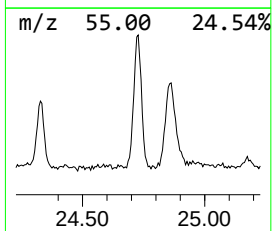
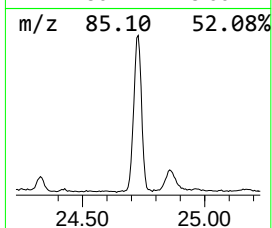
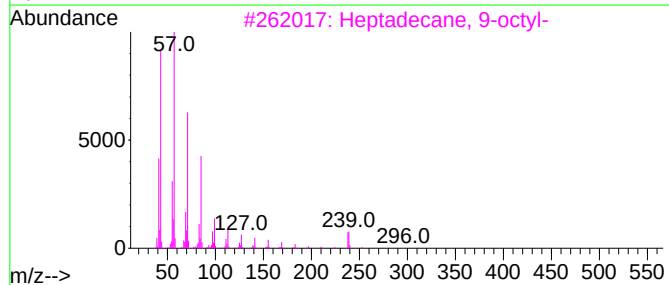
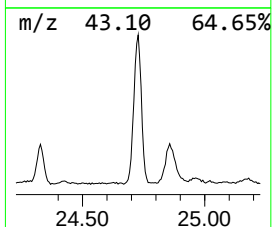
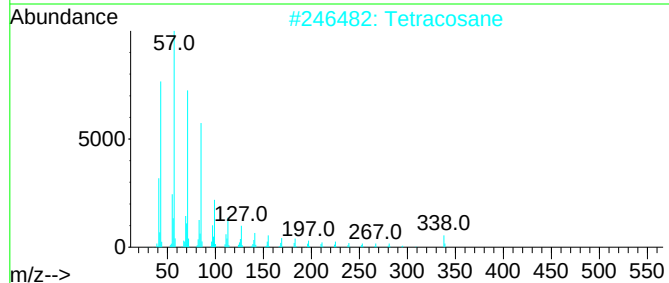
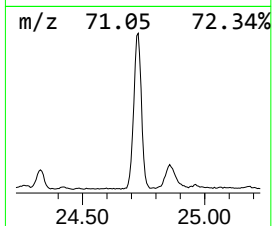
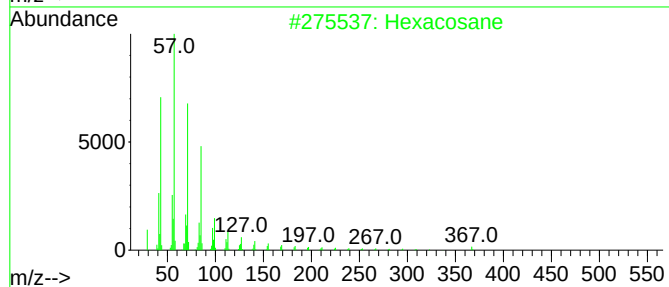
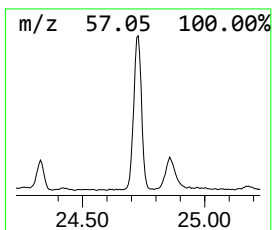
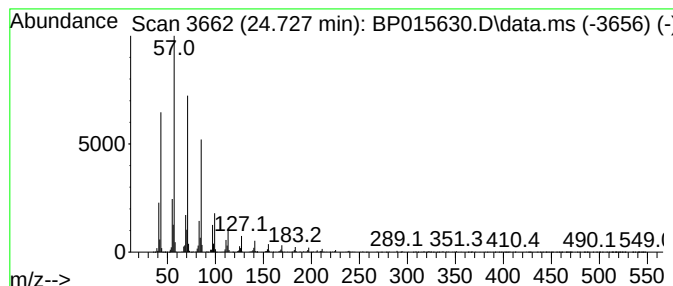
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.727	25.06 ng/ul	2133120	Perylene-d12	24.116	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexacosane	366	C26H54	000630-01-3	94
2	Tetracosane	338	C24H50	000646-31-1	94
3	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	91
4	Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5	Heptacosane	380	C27H56	000593-49-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
Acq On : 20 Jun 2023 19:52
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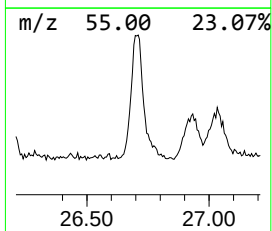
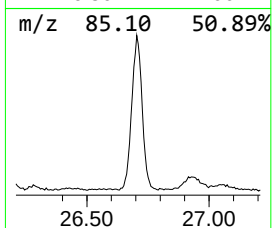
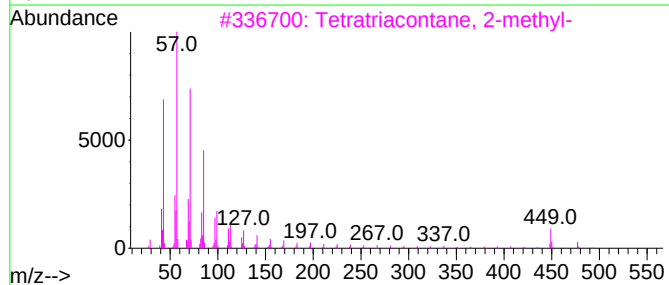
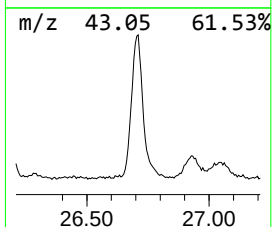
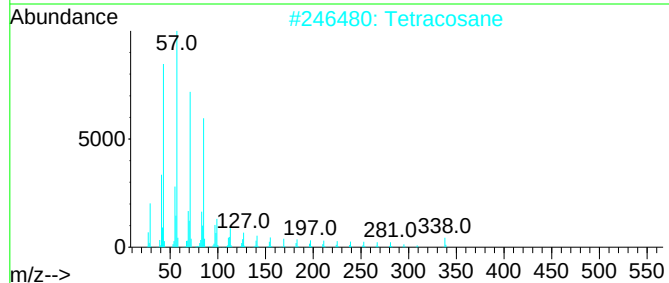
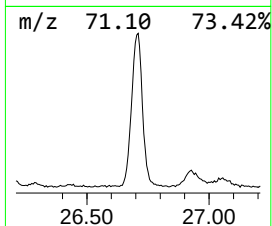
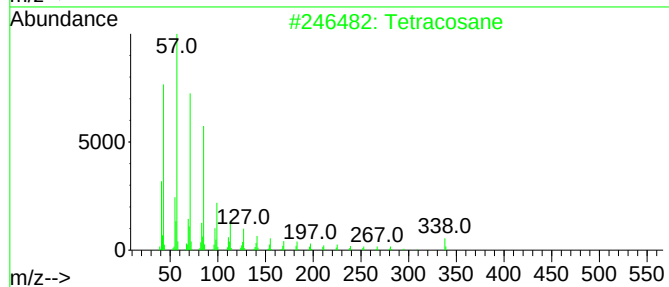
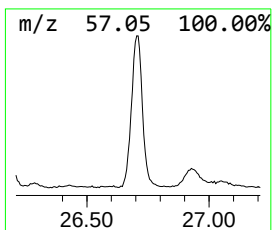
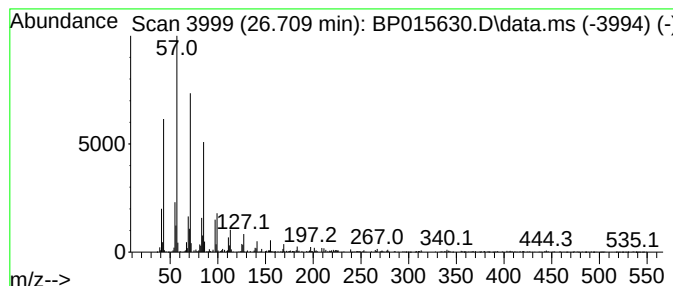
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.709	15.71 ng/ul	1337690	Perylene-d12	24.116

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetracosane	338	C24H50	000646-31-1	97
2			Tetracosane	338	C24H50	000646-31-1	94
3			Tetratriacontane, 2-methyl-	493	C35H72	014167-65-8	93
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Tetratetracontane	619	C44H90	007098-22-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062023\
Data File : BP015630.D
Acq On : 20 Jun 2023 19:52
Operator : MA/JU
Sample : 03080-19
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

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					#	RT	Resp	Conc
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Benzophenone	16.087	7.0	ng/ul	432419	3	14.728	1231400	20.0
n-Hexadecanoic ...	18.339	14.6	ng/ul	1191280	4	17.486	1627760	20.0
Octadecanoic acid	19.569	4.0	ng/ul	426466	5	21.569	2136420	20.0
(DEL) Alkane: S...	21.239	8.3	ng/ul	880977	5	21.569	2136420	20.0
(DEL) Alkane: S...	22.163	3.2	ng/ul	342611	5	21.569	2136420	20.0
(DEL) Alkane: S...	23.280	12.8	ng/ul	1086190	6	24.116	1702580	20.0
(DEL) Alkane: S...	24.727	25.1	ng/ul	2133120	6	24.116	1702580	20.0
(DEL) Alkane: S...	26.709	15.7	ng/ul	1337690	6	24.116	1702580	20.0