

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
 Data File : BP015828.D
 Acq On : 30 Jun 2023 01:25
 Operator : MA/JU
 Sample : 03161-19
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFX5

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-BP062623.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP015828.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.381	29	33	41	rVB	37493	61973	1.22%	0.087%
2	3.834	107	110	122	rBV	247826	417077	8.24%	0.584%
3	3.928	122	126	134	rVB	171209	246559	4.87%	0.345%
4	4.005	136	139	142	rBV	17372	19545	0.39%	0.027%
5	4.046	142	146	153	rVB	48127	75303	1.49%	0.106%
6	4.728	259	262	268	rVB4	24010	35599	0.70%	0.050%
7	5.099	322	325	331	rVB2	10842	15116	0.30%	0.021%
8	5.222	342	346	352	rBV3	11418	22404	0.44%	0.031%
9	6.022	477	482	488	rVB9	9665	18479	0.37%	0.026%
10	6.699	590	597	604	rVB8	13955	32770	0.65%	0.046%
11	7.010	645	650	654	rVB6	11030	19046	0.38%	0.027%
12	7.122	663	669	671	rBV6	8463	15701	0.31%	0.022%
13	7.240	682	689	707	rBV	327530	900852	17.80%	1.262%
14	7.381	707	713	724	rBV	396148	702255	13.88%	0.984%
15	7.587	741	748	763	rBV	501955	994446	19.65%	1.393%
16	7.810	781	786	790	rVB3	16743	27489	0.54%	0.039%
17	8.052	821	827	844	rBV	751830	1460182	28.85%	2.046%
18	8.799	945	954	973	rBV	291912	815449	16.11%	1.143%
19	9.252	1023	1031	1046	rBV	420252	954655	18.86%	1.338%
20	9.393	1053	1055	1065	rVB	15785	38004	0.75%	0.053%
21	9.587	1085	1088	1093	rVV5	10240	16050	0.32%	0.022%
22	9.975	1144	1154	1170	rBV	294831	787156	15.55%	1.103%
23	10.169	1185	1187	1197	rVB	12508	26396	0.52%	0.037%
24	10.269	1200	1204	1209	rBV8	7709	14812	0.29%	0.021%
25	10.551	1245	1252	1278	rVV	319549	1082195	21.38%	1.516%
26	10.887	1301	1309	1325	rVV	1052302	2180167	43.08%	3.055%
27	11.075	1335	1341	1370	rVB2	99389	395975	7.82%	0.555%
28	11.875	1473	1477	1483	rVB4	16086	25024	0.49%	0.035%
29	12.087	1506	1513	1518	rVB9	7107	17410	0.34%	0.024%
30	12.275	1539	1545	1554	rVV6	12592	29155	0.58%	0.041%
31	12.563	1591	1594	1606	rVB3	14038	31014	0.61%	0.043%
32	12.775	1625	1630	1637	rBV5	12693	29715	0.59%	0.042%
33	13.210	1699	1704	1709	rBV4	15068	26978	0.53%	0.038%
34	13.375	1726	1732	1735	rBV8	7539	12407	0.25%	0.017%
35	13.422	1735	1740	1742	rBV6	12069	18331	0.36%	0.026%
36	13.451	1742	1745	1749	rVB6	17279	21732	0.43%	0.030%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
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37	13.504	1749	1754	1758	rBV2	18523	29751	0.59%	0.042%
38	13.563	1758	1764	1772	rVB3	144706	253439	5.01%	0.355%
39	13.810	1802	1806	1810	rVB7	9431	13381	0.26%	0.019%
40	13.892	1812	1820	1826	rBV7	10163	30352	0.60%	0.043%
41	14.010	1835	1840	1849	rBV5	62167	108854	2.15%	0.153%
42	14.116	1849	1858	1870	rBV	1088122	1853751	36.63%	2.598%
43	14.234	1875	1878	1883	rVV5	40840	61922	1.22%	0.087%
44	14.292	1883	1888	1893	rVB9	12726	23020	0.45%	0.032%
45	14.398	1900	1906	1919	rBV	1349805	2304101	45.53%	3.229%
46	14.522	1922	1927	1931	rVB7	29133	50510	1.00%	0.071%
47	14.569	1931	1935	1941	rVB	37442	51500	1.02%	0.072%
48	14.645	1943	1948	1951	rBV5	21549	30267	0.60%	0.042%
49	14.704	1951	1958	1979	rVV	1800348	2922137	57.74%	4.095%
50	14.934	1991	1997	2001	rBV3	31540	57811	1.14%	0.081%
51	14.992	2001	2007	2024	rVV3	152591	579158	11.44%	0.812%
52	15.116	2025	2028	2036	rVB6	37162	77849	1.54%	0.109%
53	15.475	2084	2089	2093	rBV6	14664	30070	0.59%	0.042%
54	15.522	2094	2097	2103	rVB	37970	48754	0.96%	0.068%
55	15.704	2121	2128	2145	rBV	1674972	2923975	57.77%	4.097%
56	15.869	2151	2156	2165	rBV2	213808	482604	9.54%	0.676%
57	16.069	2185	2190	2199	rBV	338189	564134	11.15%	0.791%
58	16.163	2203	2206	2211	rVB6	22226	35577	0.70%	0.050%
59	16.410	2240	2248	2252	rBV4	69224	161287	3.19%	0.226%
60	16.551	2267	2272	2278	rBV6	48588	87603	1.73%	0.123%
61	16.610	2278	2282	2289	rVB3	70013	118286	2.34%	0.166%
62	16.845	2318	2322	2325	rBV3	32471	42877	0.85%	0.060%
63	17.186	2376	2380	2385	rBV2	44437	63626	1.26%	0.089%
64	17.333	2401	2405	2412	rBV2	133839	224491	4.44%	0.315%
65	17.404	2413	2417	2421	rVV2	71025	98875	1.95%	0.139%
66	17.463	2421	2427	2432	rBV	2159610	3265842	64.53%	4.576%
67	17.504	2432	2434	2440	rVV	262812	417195	8.24%	0.585%
68	17.569	2440	2445	2457	rVB2	1676495	2769931	54.73%	3.881%
69	18.210	2551	2554	2559	rVB4	73194	92813	1.83%	0.130%
70	18.263	2559	2563	2567	rVB	111430	131994	2.61%	0.185%
71	18.316	2568	2572	2580	rBV	1753292	2454966	48.51%	3.440%
72	18.504	2599	2604	2610	rBV3	245300	428312	8.46%	0.600%
73	18.604	2616	2621	2625	rBV3	112632	204119	4.03%	0.286%
74	19.192	2718	2721	2727	rBV2	131100	228259	4.51%	0.320%
75	19.310	2734	2741	2746	rBV4	434778	868421	17.16%	1.217%
76	19.404	2753	2757	2759	rBV4	246316	323722	6.40%	0.454%

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77	19.433	2759	2762	2763	rVV	238409	289481	5.72%	0.406%
78	19.463	2763	2767	2773	rVV	786250	1360673	26.89%	1.907%
79	19.539	2777	2780	2784	rVV2	764869	1052766	20.80%	1.475%
80	19.586	2784	2788	2792	rVB	678433	894540	17.67%	1.253%
81	19.751	2813	2816	2817	rBV2	178841	195867	3.87%	0.274%
82	19.792	2818	2823	2827	rBV	2384219	3369130	66.57%	4.721%
83	19.916	2841	2844	2850	rVB3	279720	383215	7.57%	0.537%
84	20.022	2858	2862	2867	rVB	452207	539389	10.66%	0.756%
85	20.269	2900	2904	2907	rBV2	543989	687793	13.59%	0.964%
86	20.327	2911	2914	2919	rVB	414216	488132	9.64%	0.684%
87	20.539	2947	2950	2955	rBV	325477	353468	6.98%	0.495%
88	20.598	2956	2960	2963	rBV3	266264	393065	7.77%	0.551%
89	20.627	2963	2965	2970	rVB3	180364	240574	4.75%	0.337%
90	20.751	2984	2986	2989	rBV	192002	198285	3.92%	0.278%
91	20.851	3001	3003	3010	rVB	299736	367112	7.25%	0.514%
92	21.222	3060	3066	3072	rBV3	1453880	2402422	47.47%	3.366%
93	21.551	3118	3122	3127	rVB2	1982754	2757875	54.49%	3.865%
94	22.133	3218	3221	3224	rVV	606822	702887	13.89%	0.985%
95	22.174	3224	3228	3237	rVB2	1175305	2187655	43.23%	3.066%
96	23.239	3403	3409	3415	rVB	3004602	5061071	100.00%	7.092%
97	23.321	3418	3423	3434	rVB	906195	2125747	42.00%	2.979%
98	24.098	3550	3555	3571	rVB	1234446	2831679	55.95%	3.968%
99	24.674	3647	3653	3665	rVB	1739901	3819736	75.47%	5.352%
100	26.121	3894	3899	3910	rVB	1039266	2558152	50.55%	3.585%

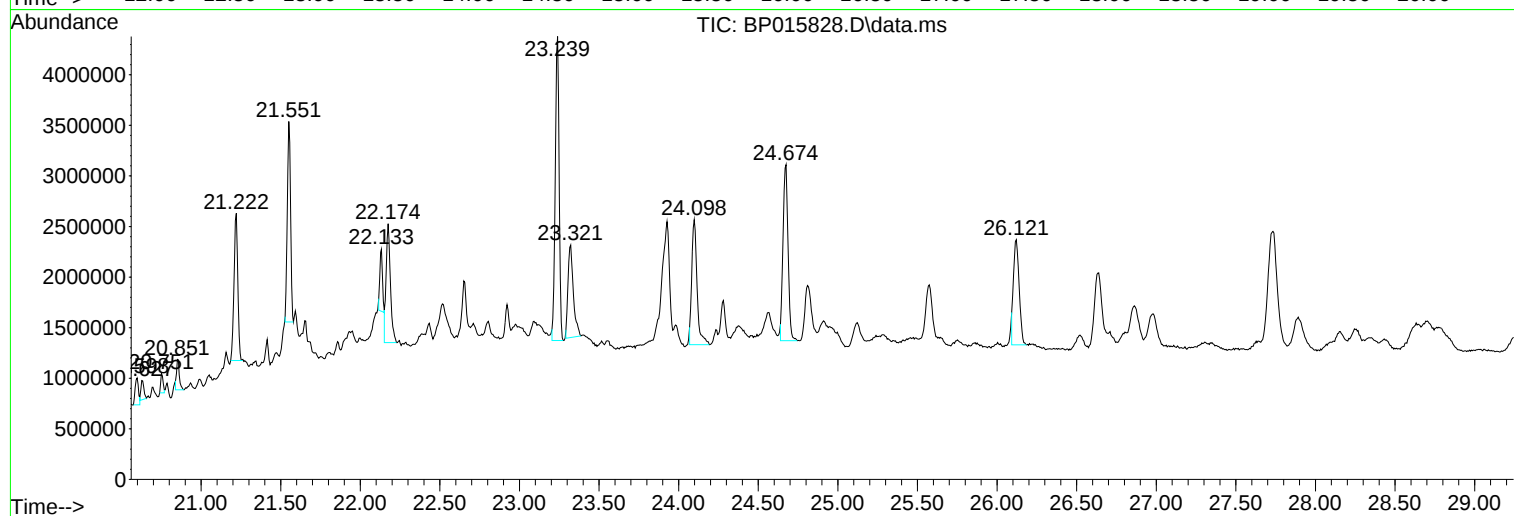
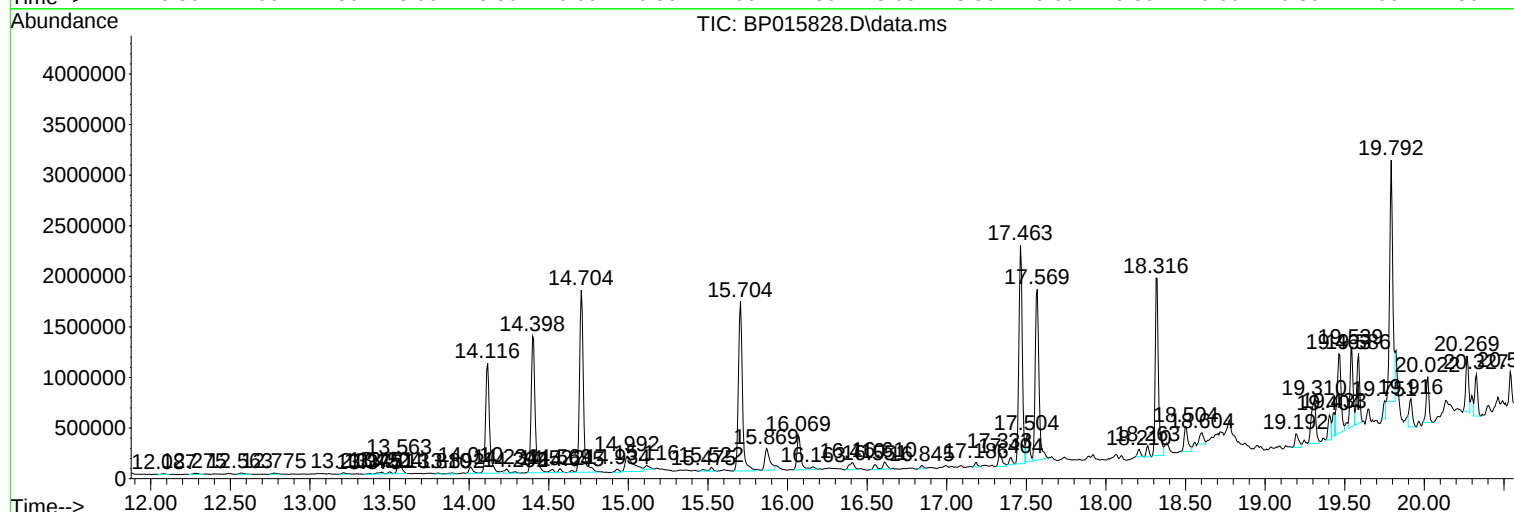
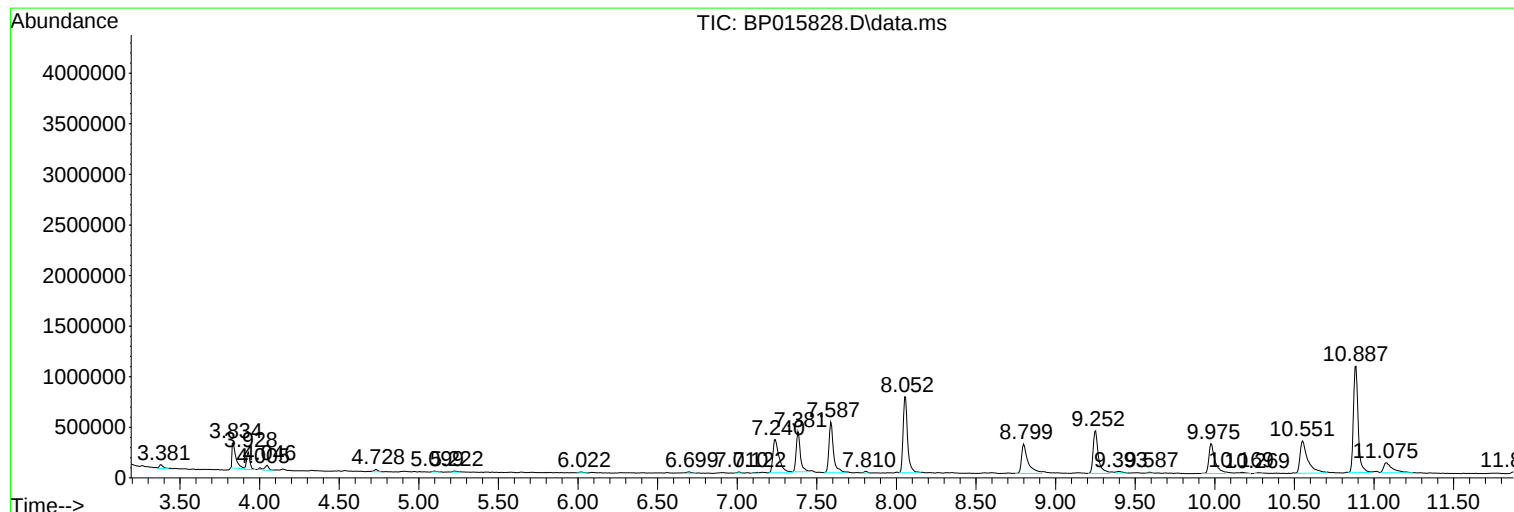
Sum of corrected areas: 71363671

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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



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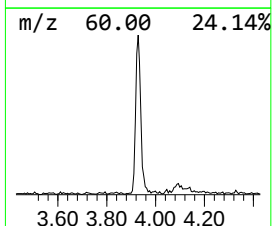
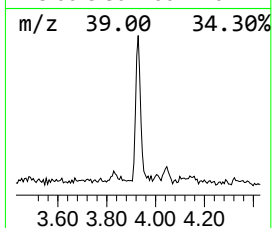
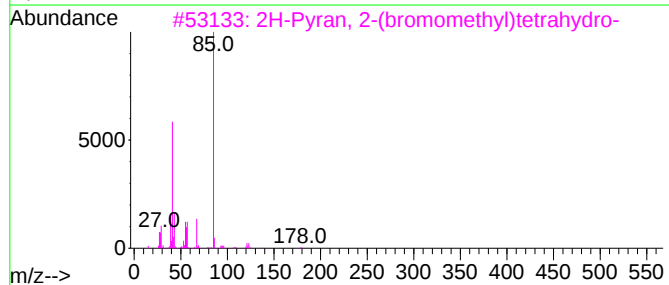
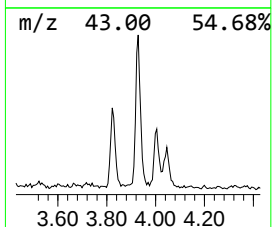
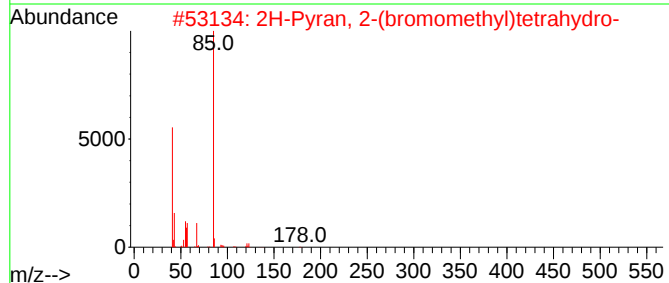
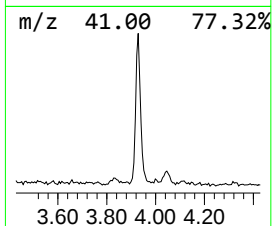
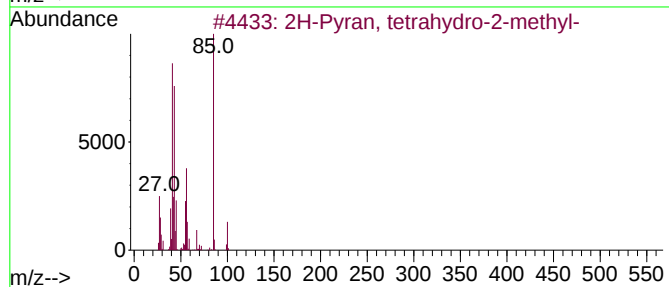
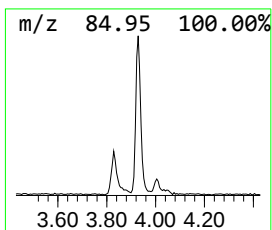
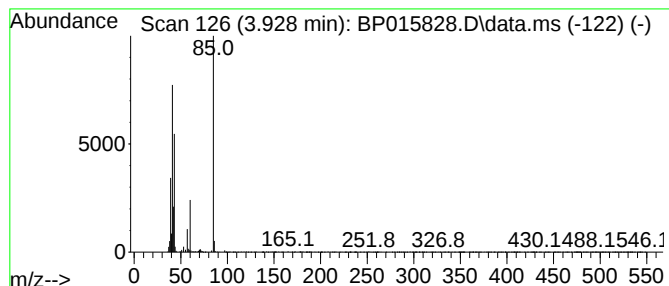
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 2H-Pyran, tetrahydro-2-methyl- Concentration Rank 15

R.T.	EstConc	Area	Relative to ISTD	R.T.
3.928	3.38 ng/ul	246559	1,4-Dichlorobenzene-d4	8.057

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	64
2			2H-Pyran, 2-(bromomethyl)tetrahy...	178	C6H11BrO	034723-82-5	42
3			2H-Pyran, 2-(bromomethyl)tetrahy...	178	C6H11BrO	034723-82-5	40
4			2H-Pyran, tetrahydro-2-methyl-	100	C6H12O	010141-72-7	39
5			Methacrylamide	85	C4H7NO	000079-39-0	36



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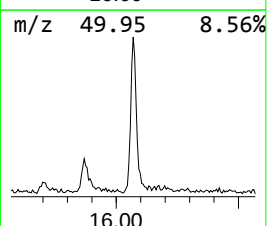
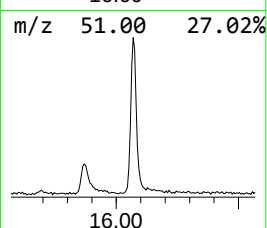
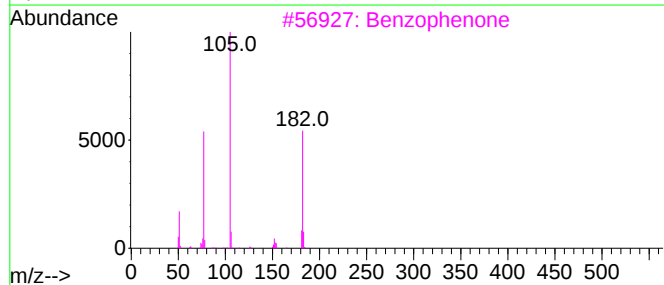
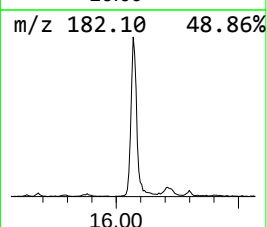
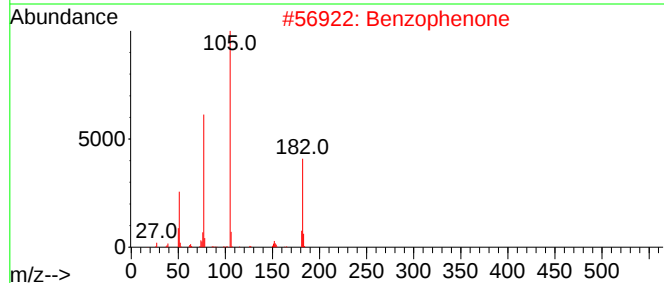
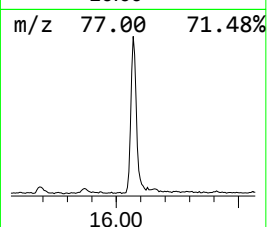
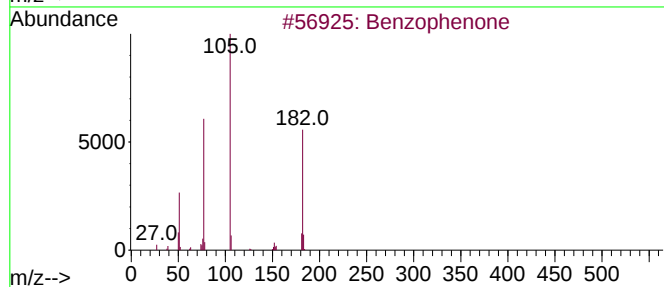
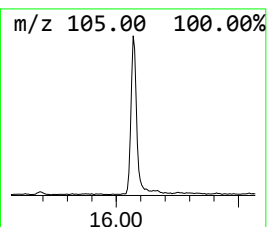
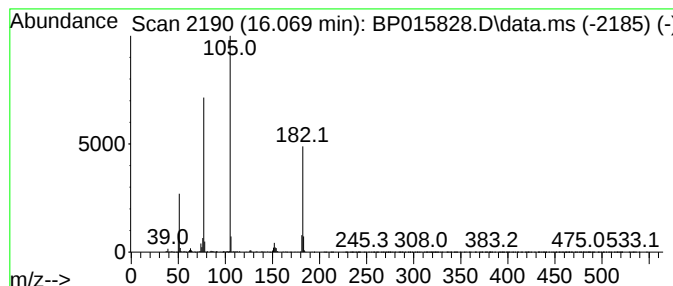
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.069	3.86 ng/ul	564134	Acenaphthene-d10	14.704

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	91



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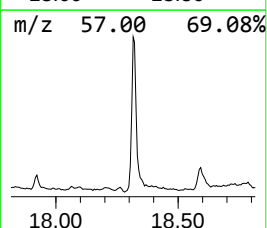
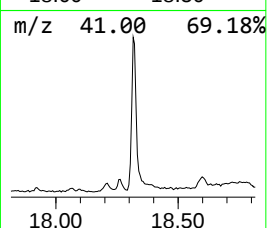
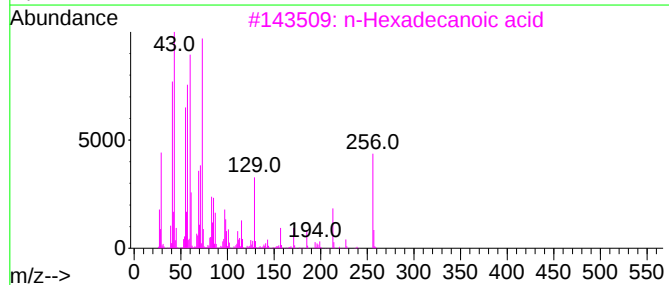
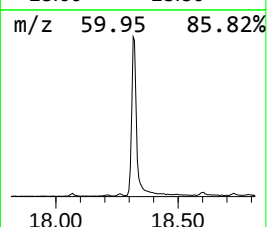
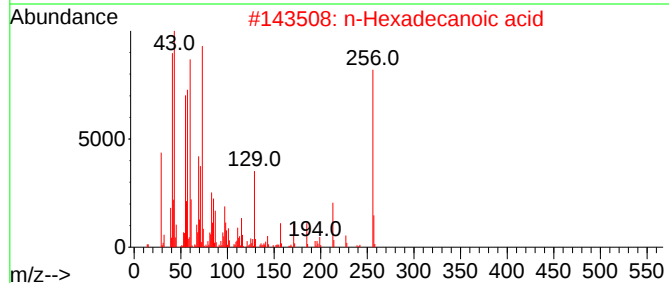
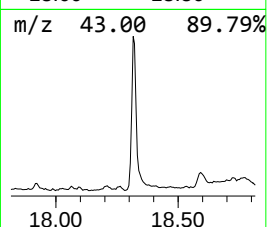
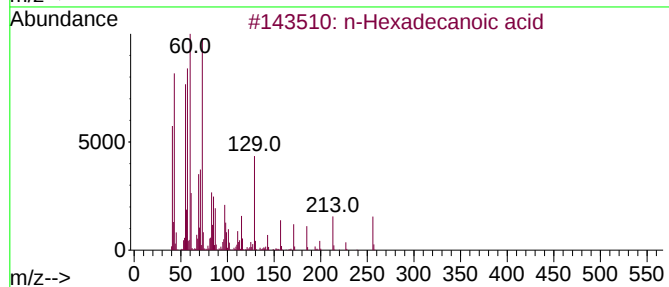
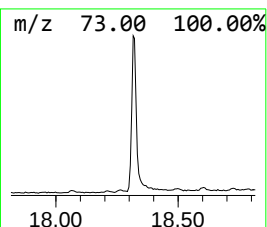
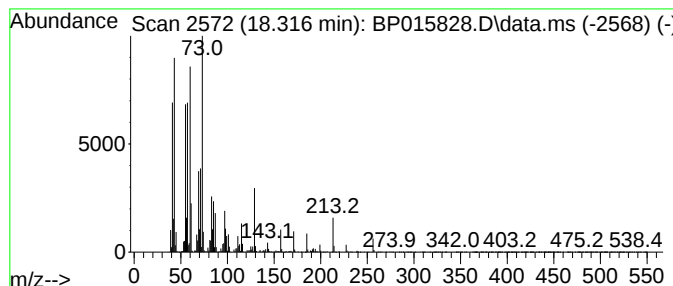
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.316	15.03 ng/ul	2454970	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	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tridecanoic acid	214	C13H26O2	000638-53-9	95
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX5

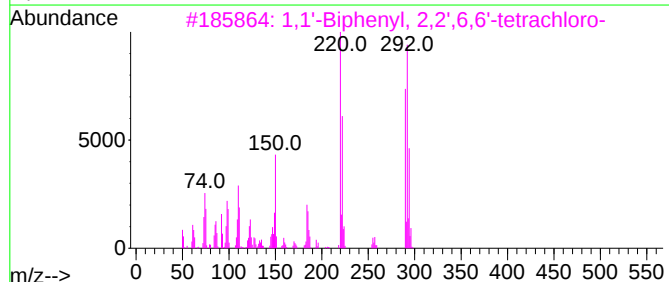
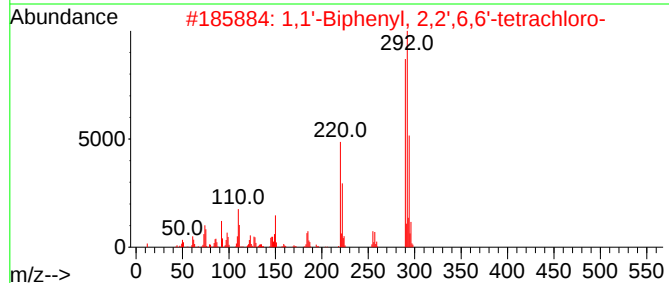
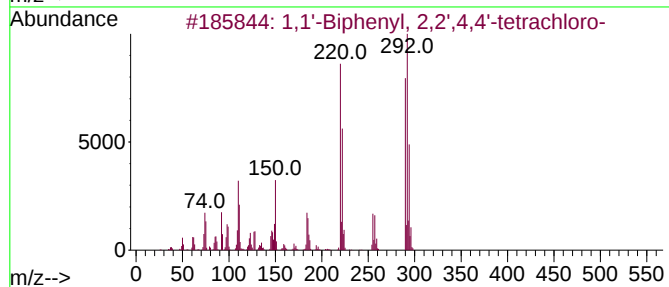
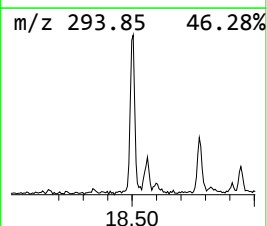
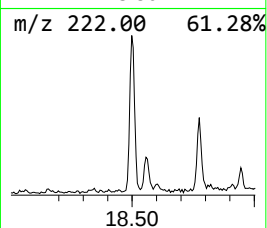
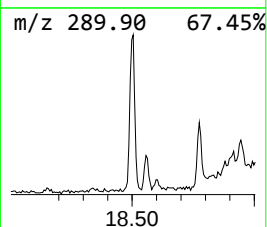
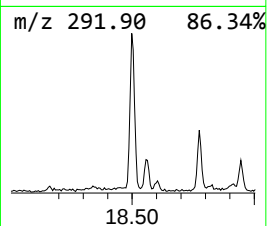
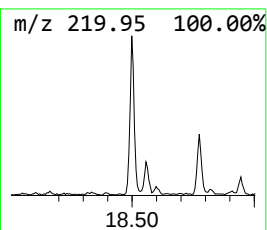
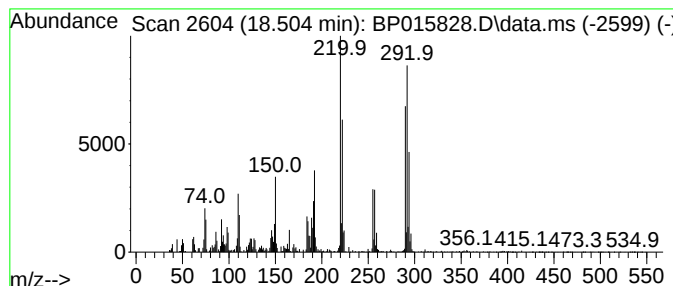
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,1'-Biphenyl, 2,2',4,4'-te... Concentration Rank 19

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.504	2.62 ng/ul	428312	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
2	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99		
3	1,1'-Biphenyl, 2,2',6,6'-tetrach...	290	C12H6Cl4	015968-05-5	99		
4	1,1'-Biphenyl, 2,2',4,4'-tetrach...	290	C12H6Cl4	002437-79-8	99		
5	1,1'-Biphenyl, 2,2',3,5'-tetrach...	290	C12H6Cl4	041464-39-5	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX5

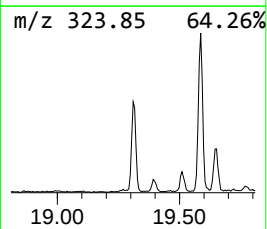
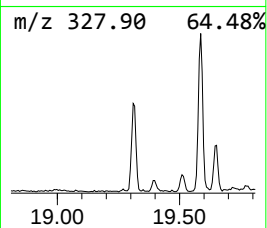
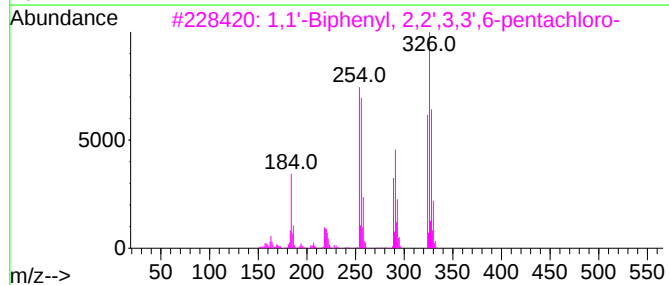
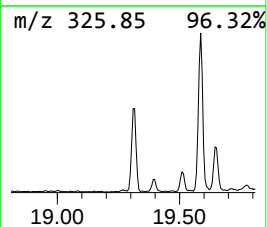
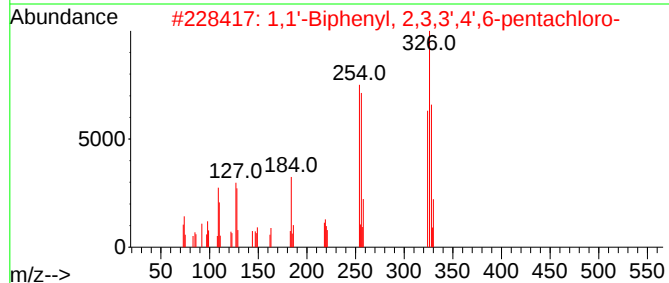
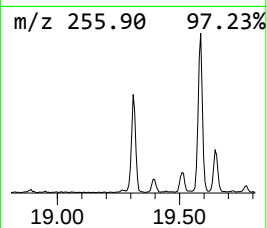
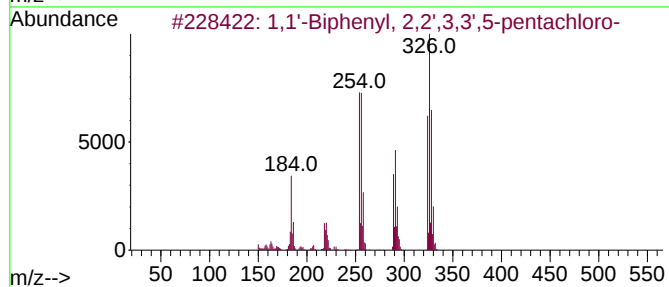
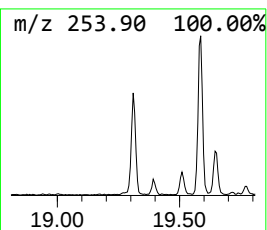
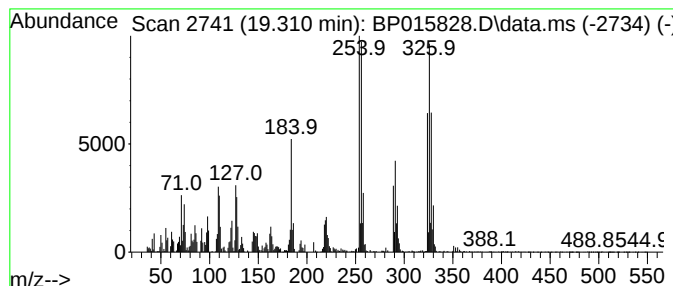
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 1,1'-Biphenyl, 2,2',3,3',5-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.310	5.32 ng/ul	868421	Phenanthrene-d10	17.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,3',5-penta...	324	C12H5Cl5	060145-20-2	99		
2	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
3	1,1'-Biphenyl, 2,2',3,3',6-penta...	324	C12H5Cl5	052663-60-2	99		
4	2,2',3,4',5-Pentachloro-1,1'-bip...	324	C12H5Cl5	068194-07-0	99		
5	1,1'-Biphenyl, 2,2',4,5,5'-penta...	324	C12H5Cl5	037680-73-2	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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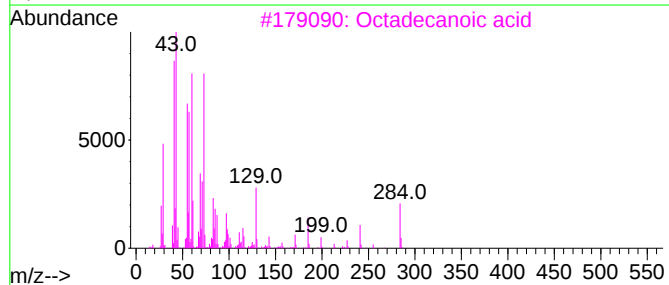
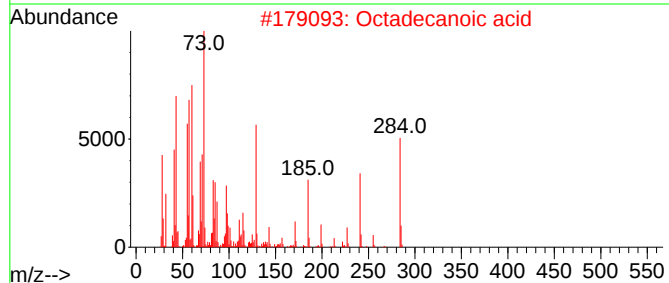
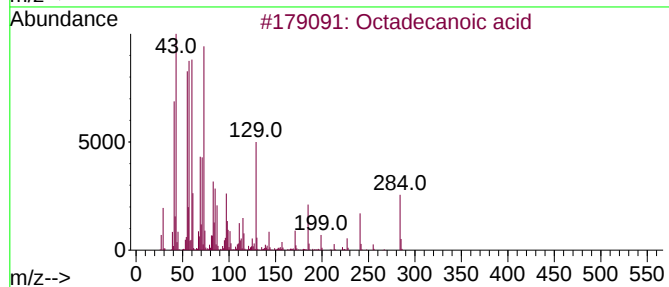
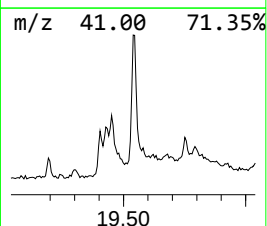
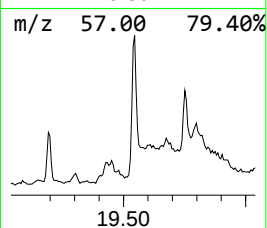
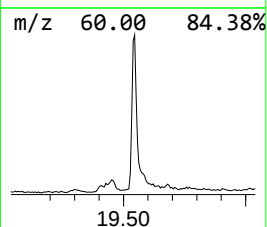
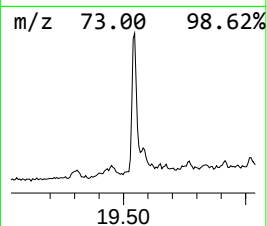
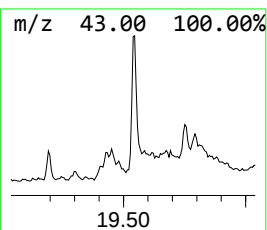
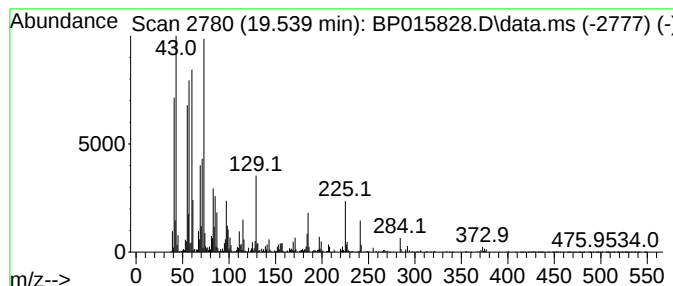
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Octadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.539	7.63 ng/ul	1052770	Chrysene-d12	21.551

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	94
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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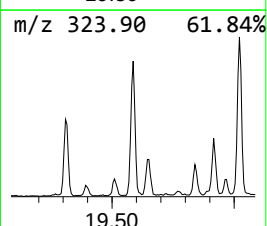
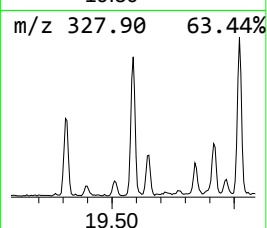
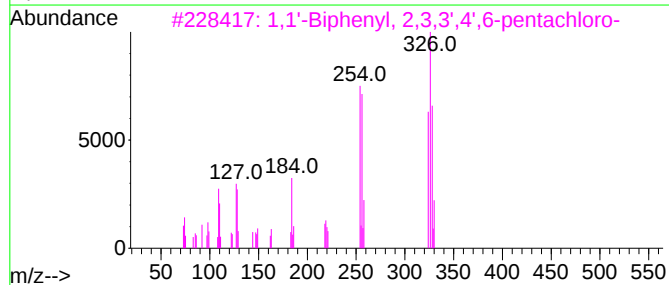
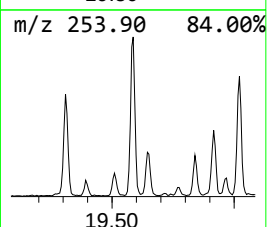
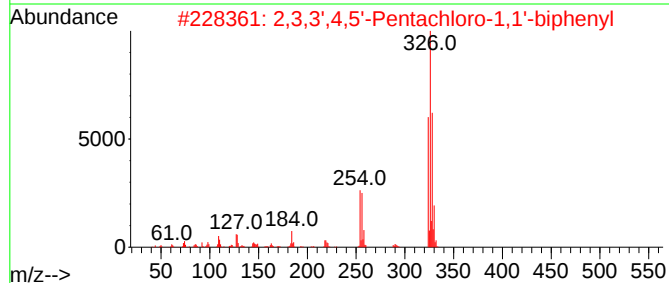
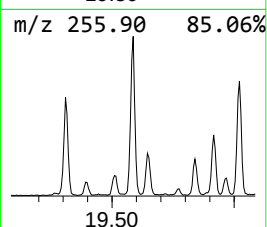
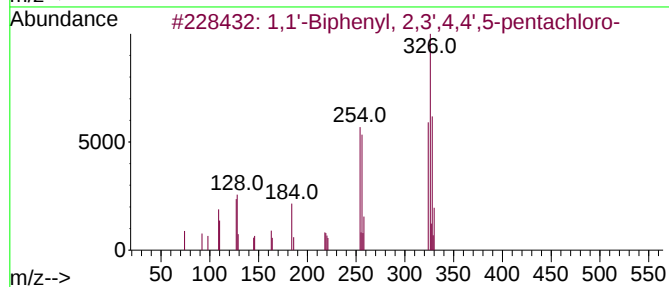
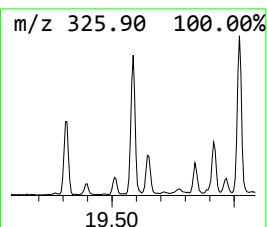
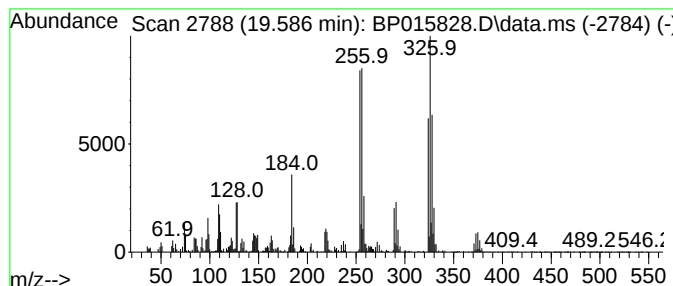
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1,1'-Biphenyl, 2,3',4,4',5-... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.586	6.49 ng/ul	894540	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99		
2	2,3,3',4,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	070362-41-3	99		
3	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
4	1,1'-Biphenyl, 2,2',4,6,6'-Penta...	324	C12H5Cl5	056558-16-8	99		
5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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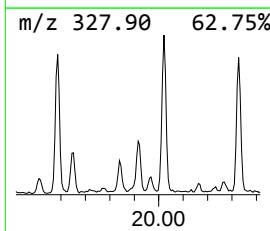
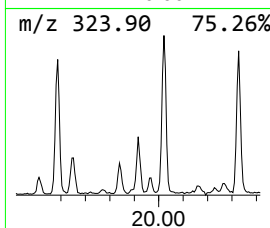
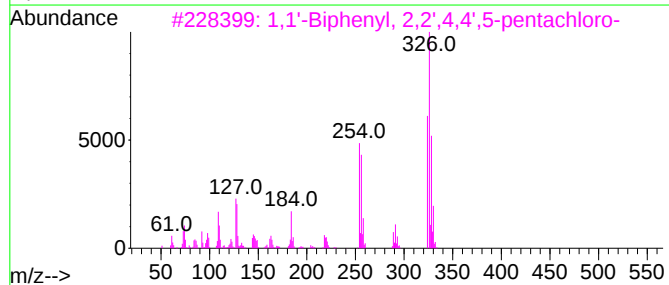
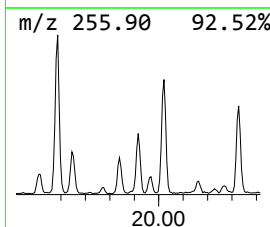
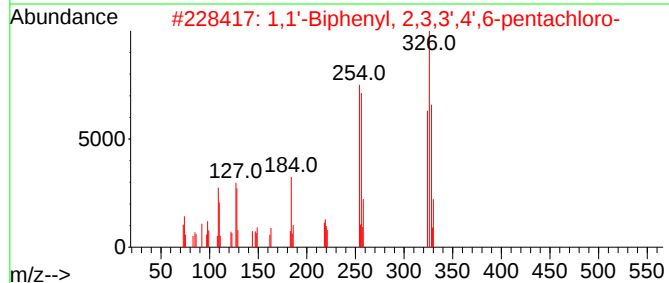
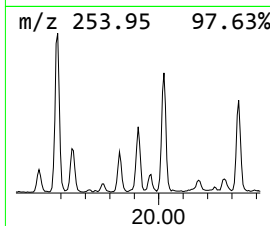
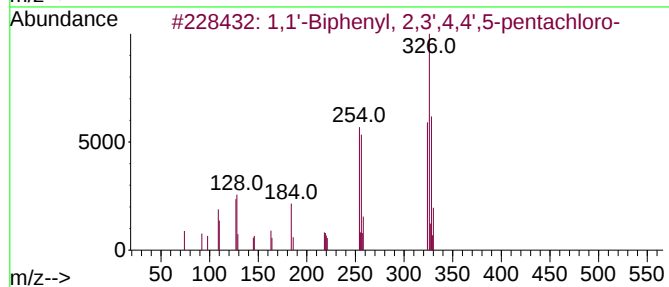
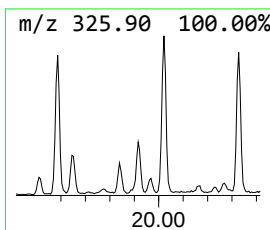
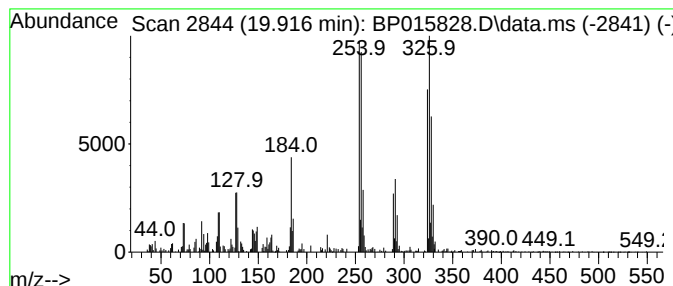
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1,1'-Biphenyl, 2,3,3',4',6-... Concentration Rank 17

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.916	2.78 ng/ul	383215	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,3',4,4',5-penta...	324	C12H5Cl5	031508-00-6	99		
2	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99		
3	1,1'-Biphenyl, 2,2',4,4',5-penta...	324	C12H5Cl5	038380-01-7	99		
4	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99		
5	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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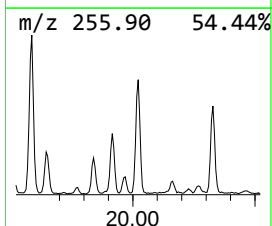
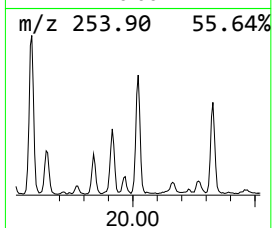
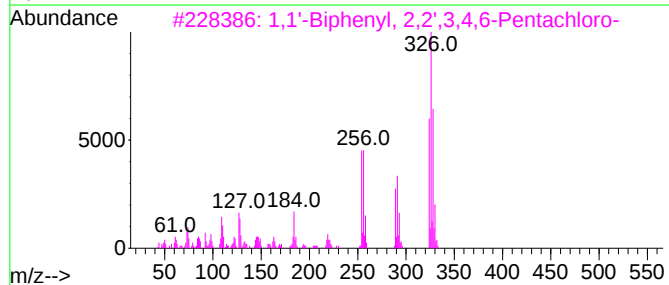
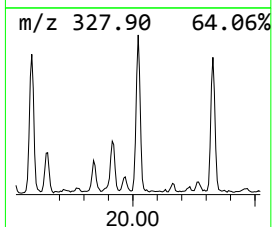
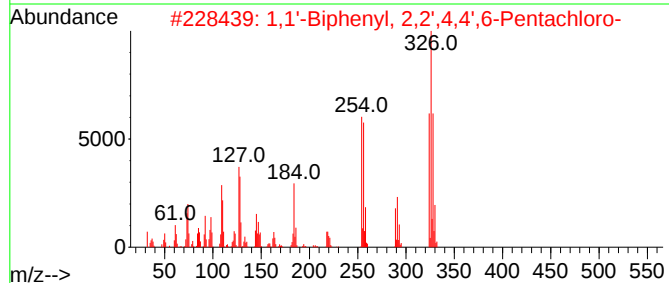
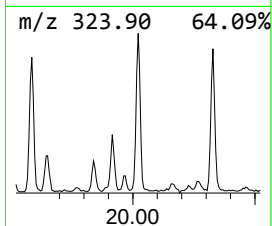
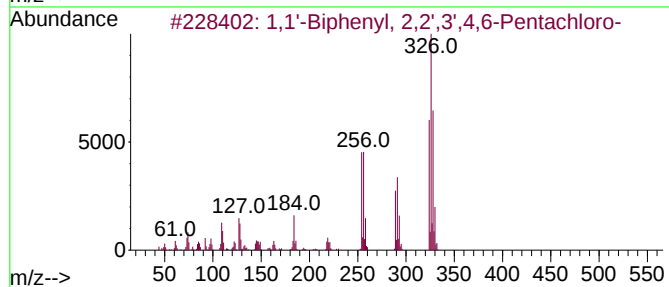
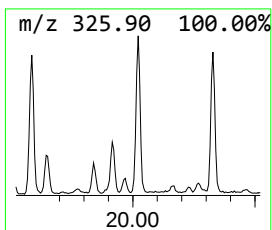
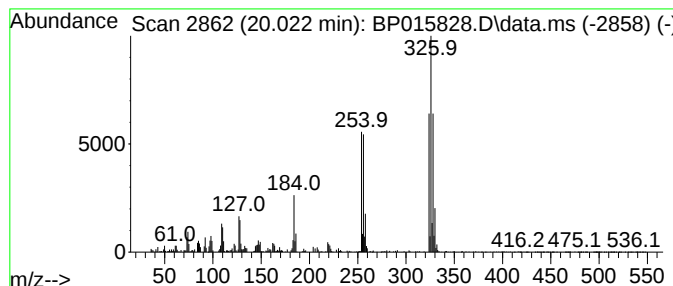
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 1,1'-Biphenyl, 2,2',3',4,6-... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.022	3.91 ng/ul	539389	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3',4,6-Penta...	324	C12H5Cl5	060233-25-2	99		
2	1,1'-Biphenyl, 2,2',4,4',6-Penta...	324	C12H5Cl5	039485-83-1	98		
3	1,1'-Biphenyl, 2,2',3,4,6-Pentac...	324	C12H5Cl5	055215-17-3	98		
4	2,3,3',5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-32-0	98		
5	1,1'-Biphenyl, 2,2',3',4,5-Penta...	324	C12H5Cl5	041464-51-1	98		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

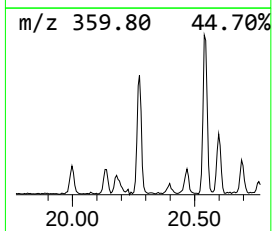
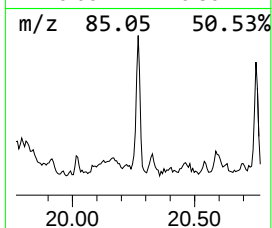
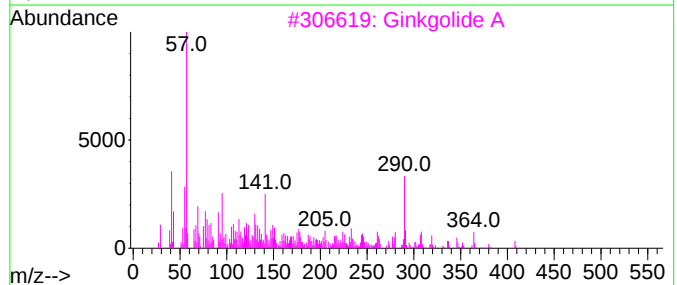
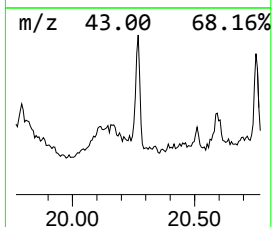
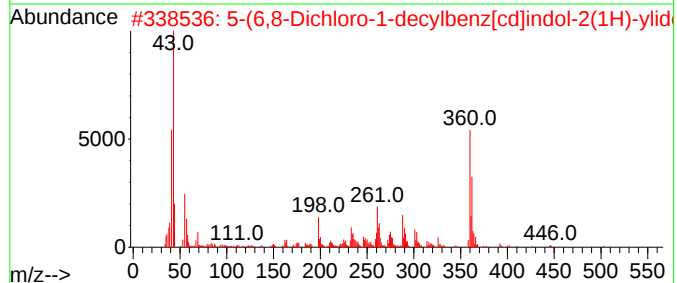
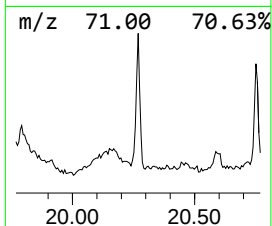
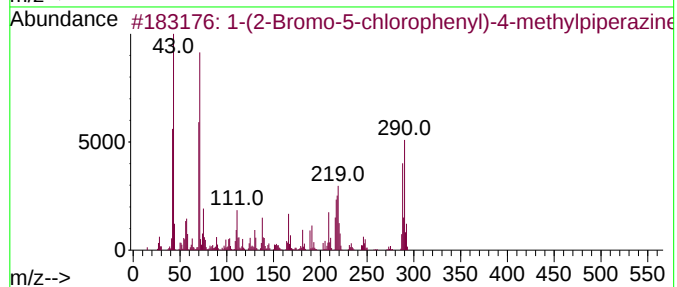
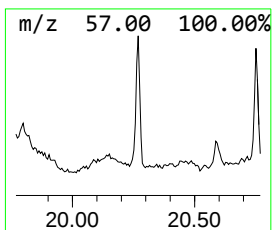
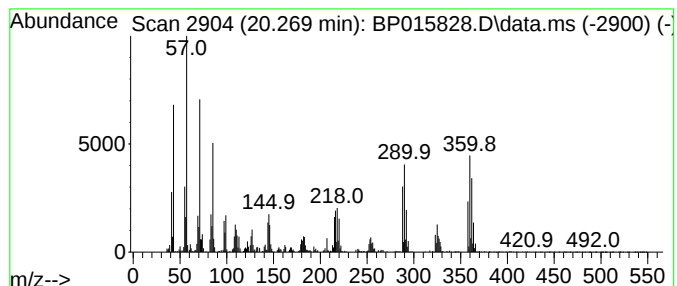
Instrument :
BNA_P
ClientSampleId :
DCFX5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-01 Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.269	4.99 ng/ul	687793	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-(2-Bromo-5-chlorophenyl)-4-met...	288	C11H14BrClN2	1261940-25-3	20
2	5-(6,8-Dichloro-1-decylbenz[cd]i...	503	C27H31Cl2NO4	303099-58-3	10
3	Ginkgolide A	408	C20H24O9	015291-75-5	9
4	Phenol, 4-nitro-2-(2-thiazolylim...	249	C10H7N3O3S	078096-09-0	9
5	1,3-Propanediol, eicosyl ethyl e...	384	C25H52O2	1000406-35-5	9



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

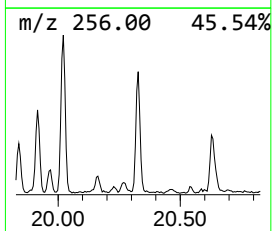
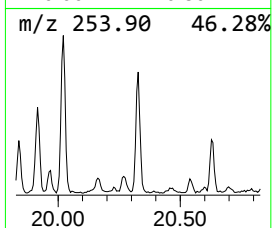
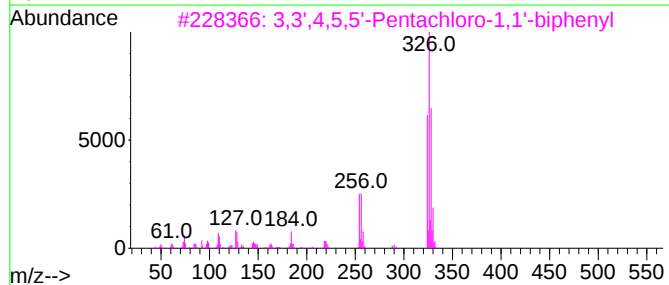
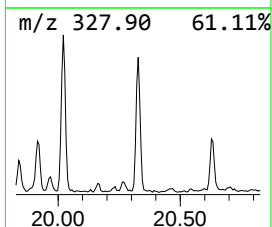
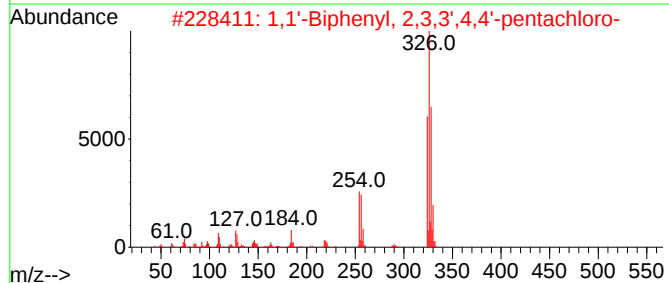
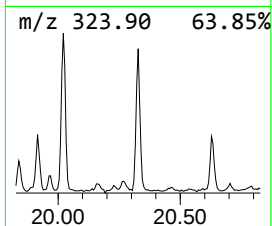
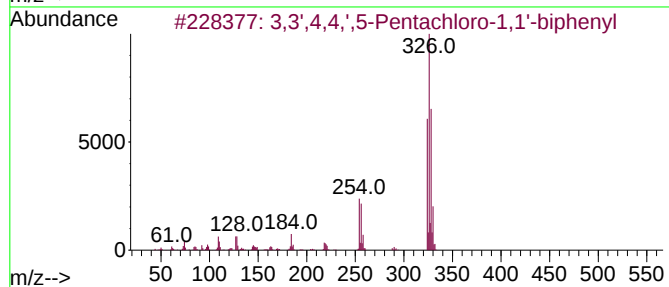
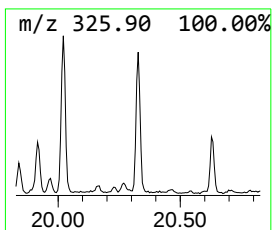
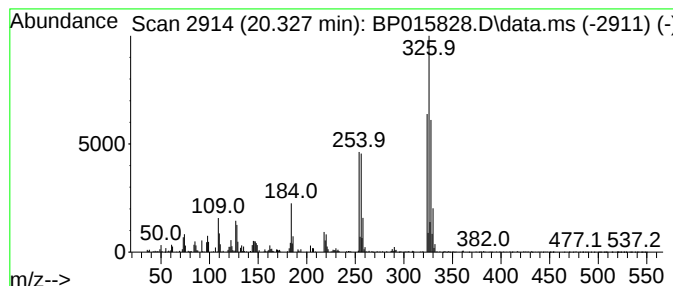
Instrument :
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 11 3,3',4,4',5-Pentachloro-1,... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.327	3.54 ng/ul	488132	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,3',4,4',5-Pentachloro-1,1'-bi...	324	C12H5Cl5	057465-28-8	99
2	1,1'-Biphenyl, 2,3,3',4,4'-penta...	324	C12H5Cl5	032598-14-4	99
3	3,3',4,5,5'-Pentachloro-1,1'-bip...	324	C12H5Cl5	039635-33-1	99
4	1,1'-Biphenyl, 2',3,4,5,5'-Penta...	324	C12H5Cl5	070424-70-3	99
5	1,1'-Biphenyl, 2,3,3',4',6-penta...	324	C12H5Cl5	038380-03-9	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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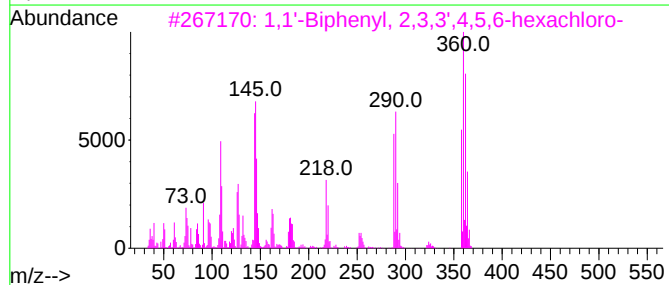
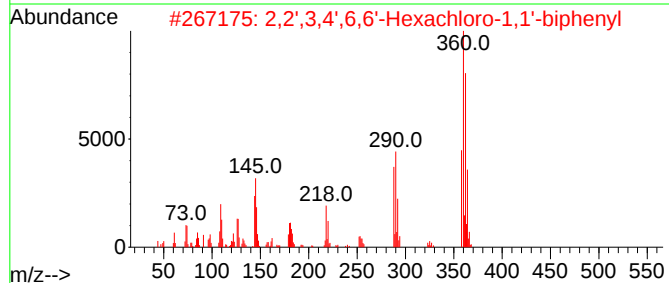
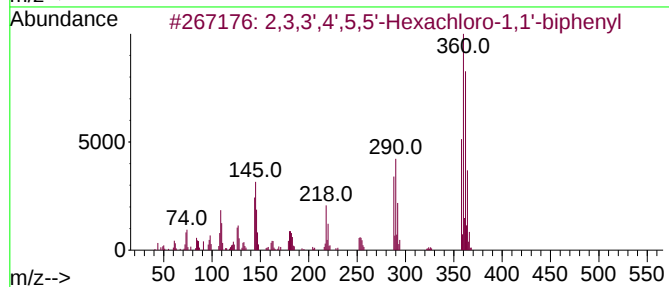
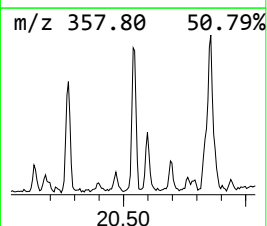
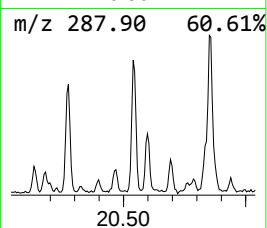
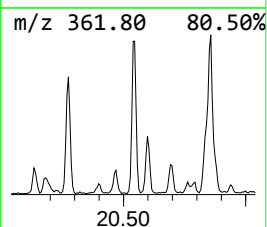
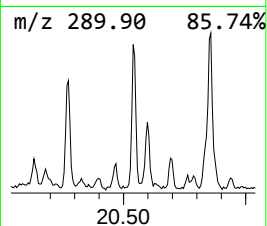
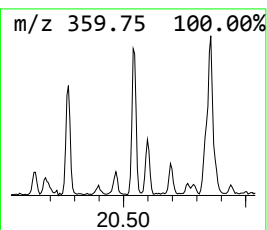
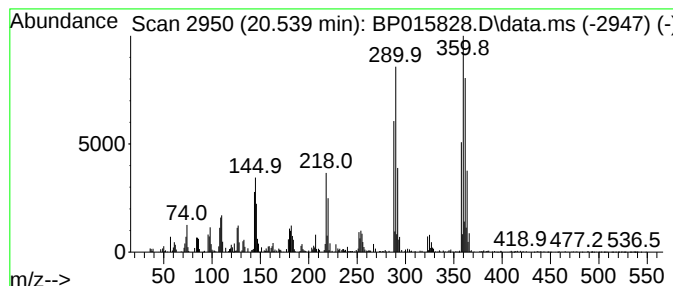
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 2,3,3',4',5,5'-Hexachloro-1... Concentration Rank 20

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.539	2.56 ng/ul	353468	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,3,3',4',5,5'-Hexachloro-1,1'-b...	358	C12H4Cl6	039635-34-2	99		
2	2,2',3,4',6,6'-Hexachloro-1,1'-b...	358	C12H4Cl6	068194-08-1	99		
3	1,1'-Biphenyl, 2,3,3',4,5,6-hexa...	358	C12H4Cl6	041411-62-5	99		
4	2,3,3',4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	074472-45-0	99		
5	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

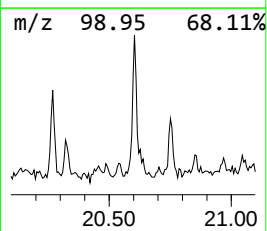
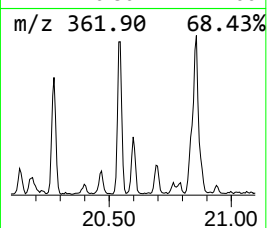
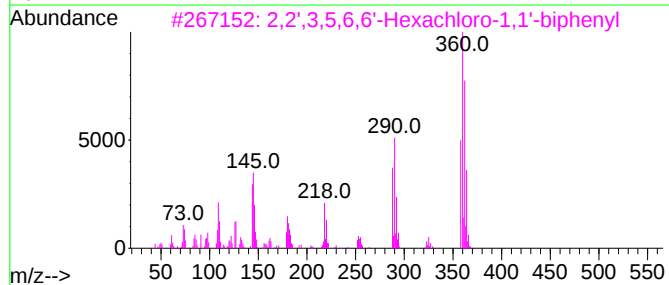
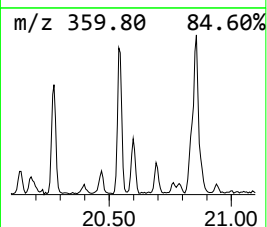
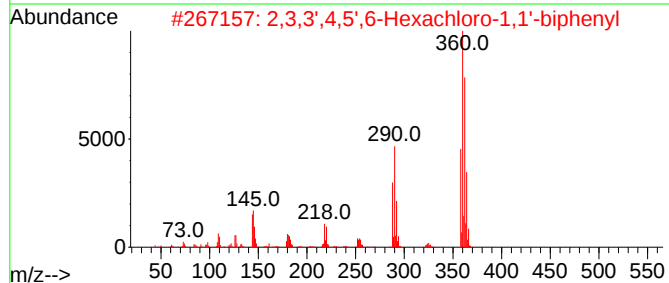
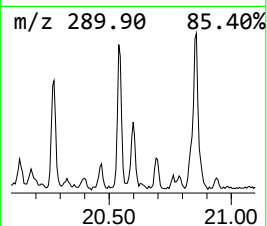
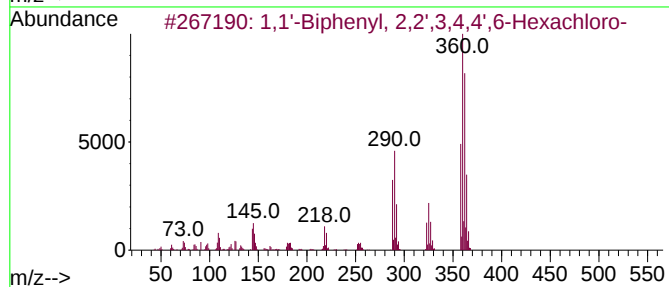
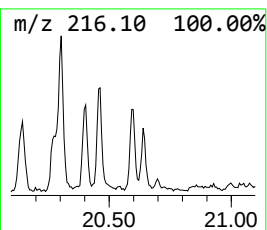
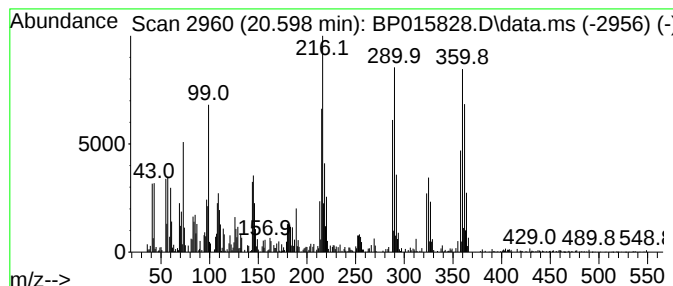
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 13 1,1'-Biphenyl, 2,2',3,4,4',... Concentration Rank 16

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.598	2.85 ng/ul	393065	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1,1'-Biphenyl, 2,2',3,4,4',6-Hex...	358	C12H4Cl6	056030-56-9	99
2	2,3,3',4,5',6-Hexachloro-1,1'-bi...	358	C12H4Cl6	074472-43-8	99
3	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99
4	1,1'-Biphenyl, 2,2',3,5,5',6-hex...	358	C12H4Cl6	052663-63-5	99
5	1,1'-Biphenyl, 2,2',3,3',4,5-hex...	358	C12H4Cl6	055215-18-4	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

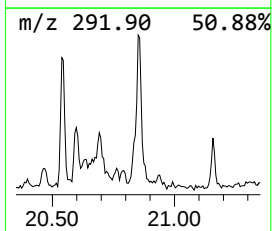
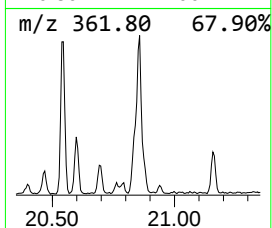
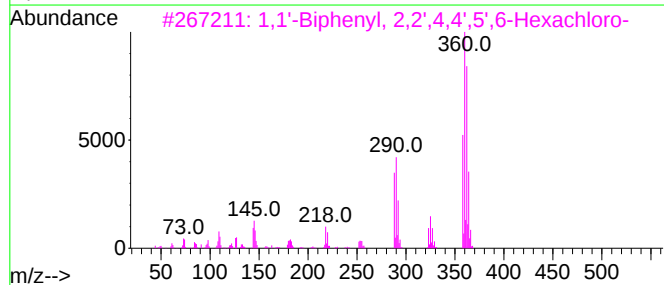
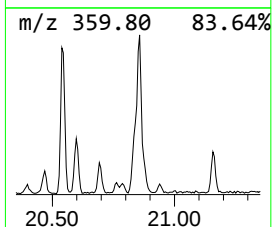
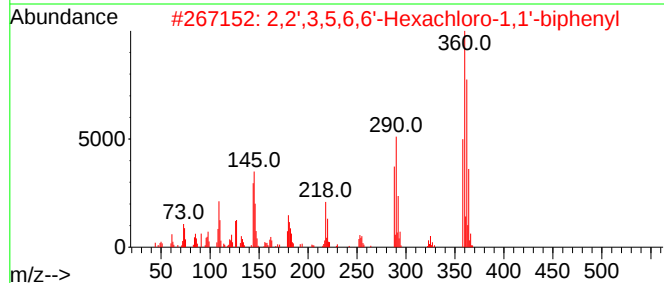
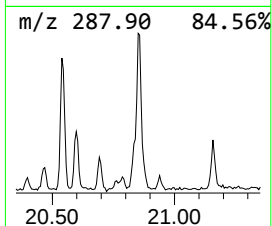
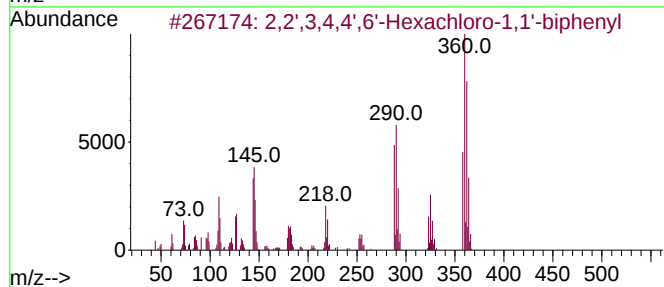
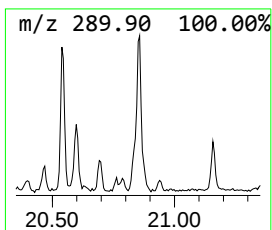
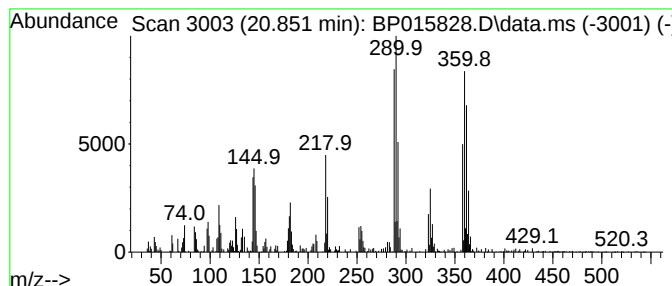
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 14 2,2',3,4,4',6'-Hexachloro-1... Concentration Rank 18

R.T.	EstConc	Area	Relative to ISTD	R.T.	
20.851	2.66 ng/ul	367112	Chrysene-d12	21.551	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2,2',3,4,4',6'-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-64-4	99
2	2,2',3,5,6,6'-Hexachloro-1,1'-bi...	358	C12H4Cl6	068194-09-2	99
3	1,1'-Biphenyl, 2,2',4,4',5',6-He...	358	C12H4Cl6	060145-22-4	99
4	1,1'-Biphenyl, 2,3,3',4,4',5-hex...	358	C12H4Cl6	038380-08-4	99
5	2,3',4,4',5',6-Hexachloro-1,1'-b...	358	C12H4Cl6	059291-65-5	99



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
DCFX5

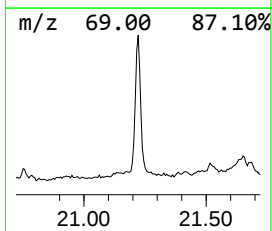
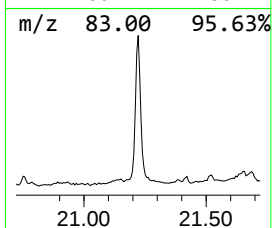
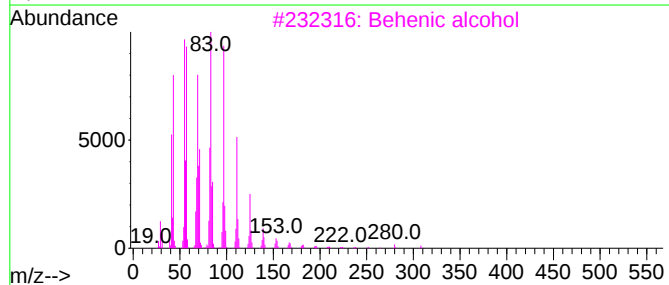
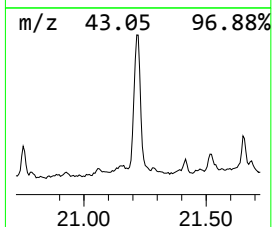
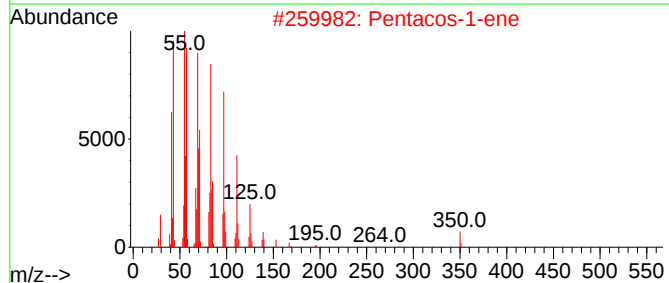
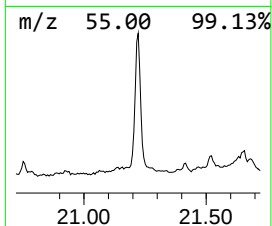
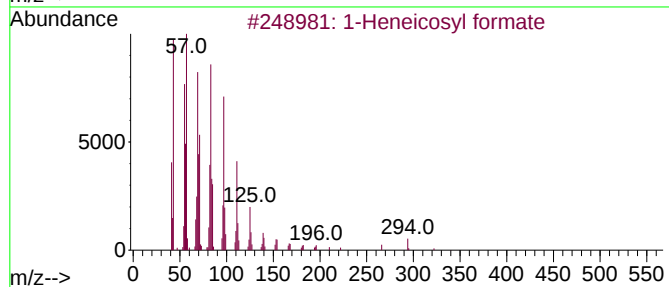
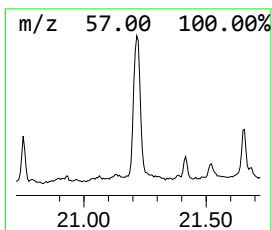
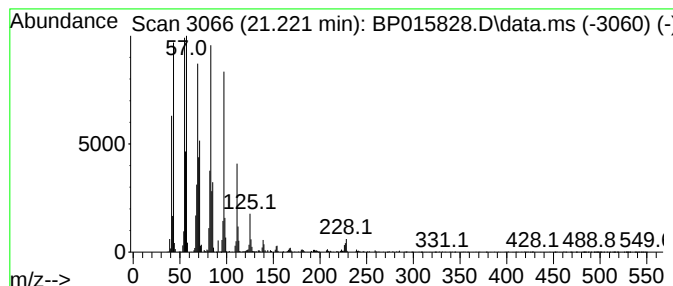
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 1-Heneicosyl formate Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.221	17.42 ng/ul	2402420	Chrysene-d12	21.551

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Heneicosyl formate	340	C22H44O2	077899-03-7	96
2			Pentacos-1-ene	350	C25H50	016980-85-1	94
3			Behenic alcohol	326	C22H46O	000661-19-8	94
4			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94
5			1-Nonadecene	266	C19H38	018435-45-5	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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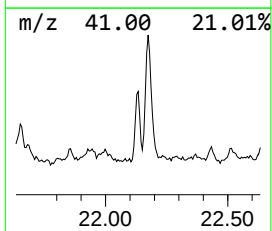
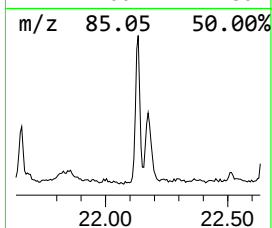
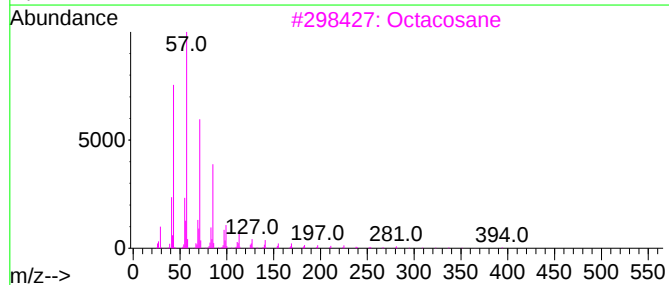
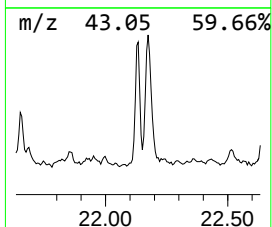
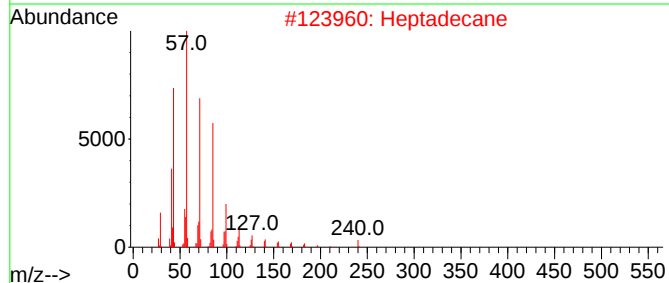
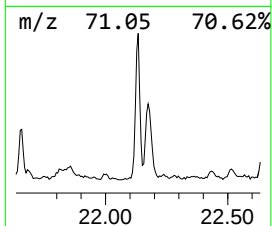
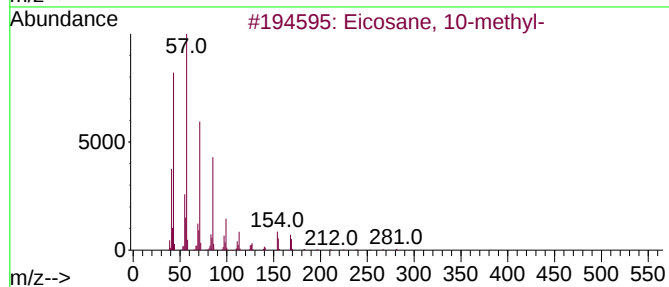
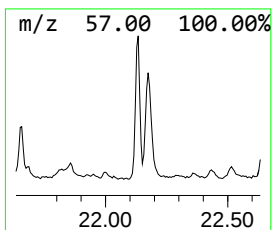
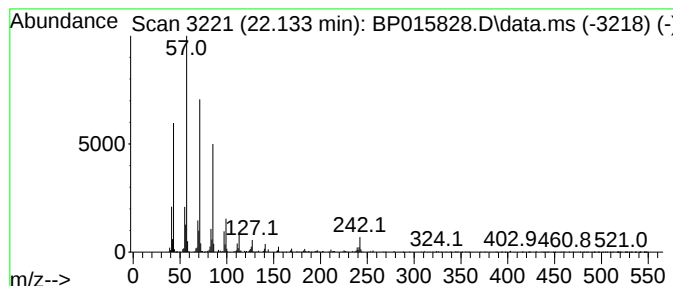
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Quant Title : SVOA CALIBRATION

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

Peak Number 16 (DEL) Alkane: Straight-Chai... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.133	5.10 ng/ul	702887	Chrysene-d12	21.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Eicosane, 10-methyl-	296	C21H44	054833-23-7	93
2		Heptadecane	240	C17H36	000629-78-7	93
3		Octacosane	394	C28H58	000630-02-4	90
4		Tetracosane	338	C24H50	000646-31-1	87
5		Heneicosane	296	C21H44	000629-94-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX5

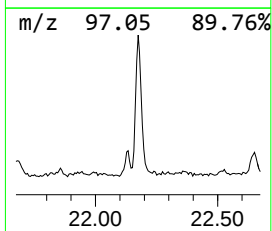
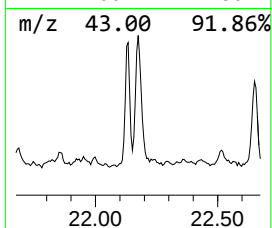
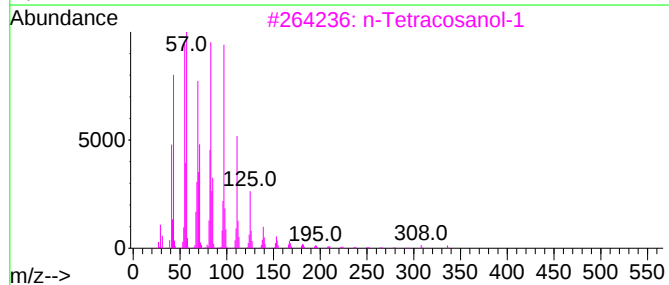
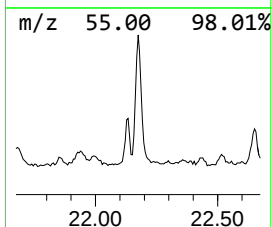
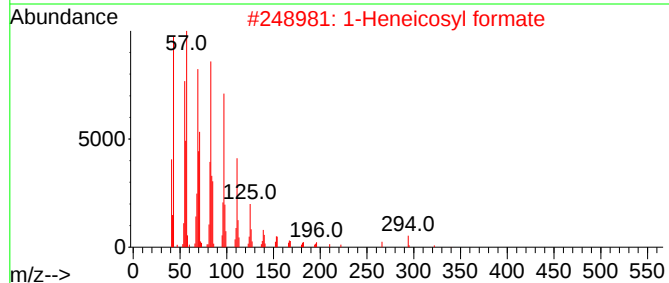
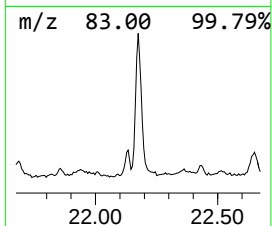
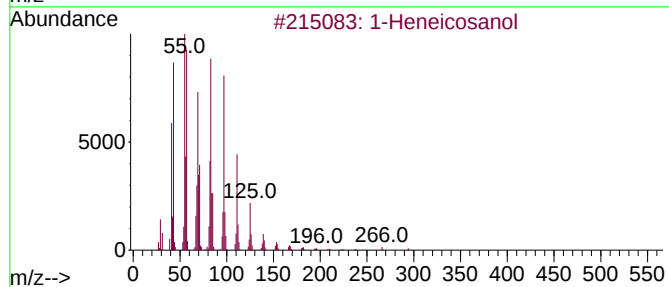
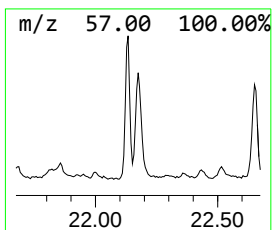
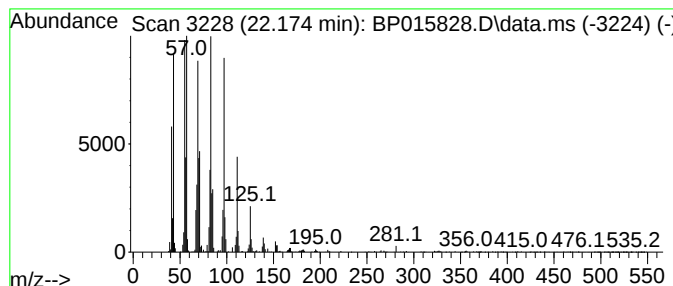
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Quant Title : SVOA CALIBRATION

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

Peak Number 17 1-Heneicosanol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.174	15.86 ng/ul	2187660	Chrysene-d12	21.551

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1-Heneicosanol	312	C21H44O	015594-90-8	95
2		1-Heneicosyl formate	340	C22H44O2	077899-03-7	95
3		n-Tetracosanol-1	354	C24H50O	000506-51-4	94
4		Behenic alcohol	326	C22H46O	000661-19-8	94
5		1-Heptacosanol	396	C27H56O	002004-39-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX5

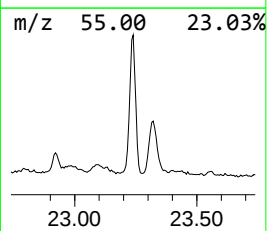
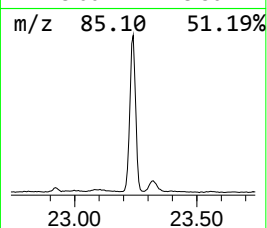
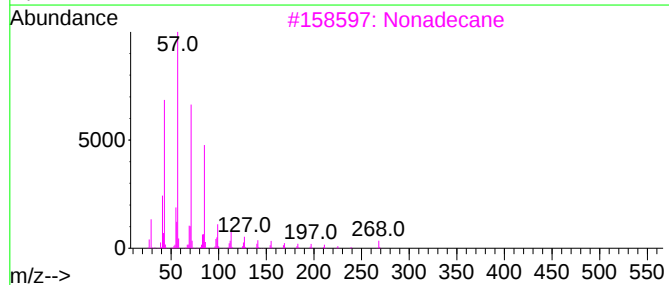
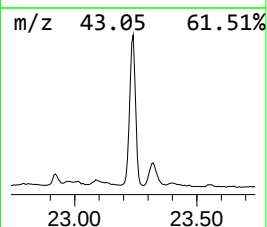
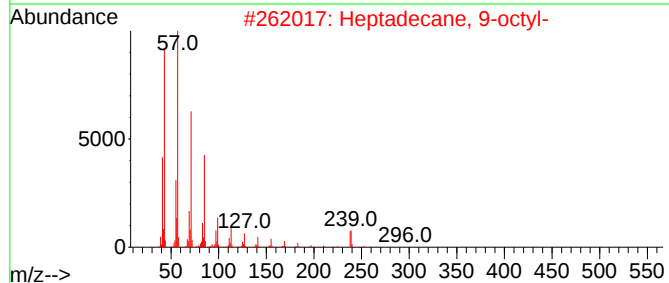
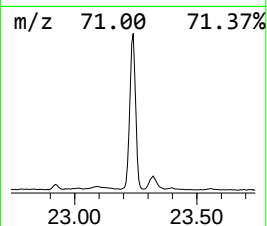
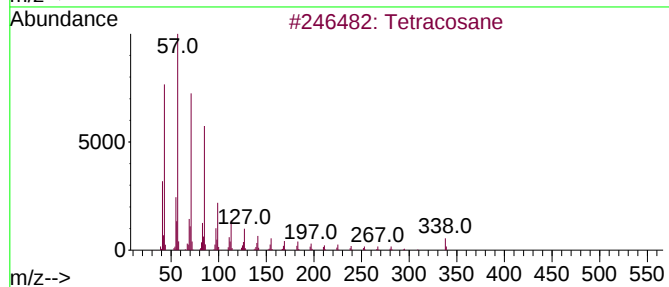
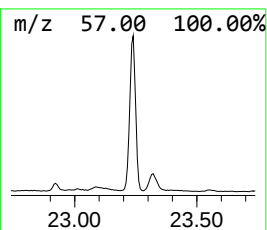
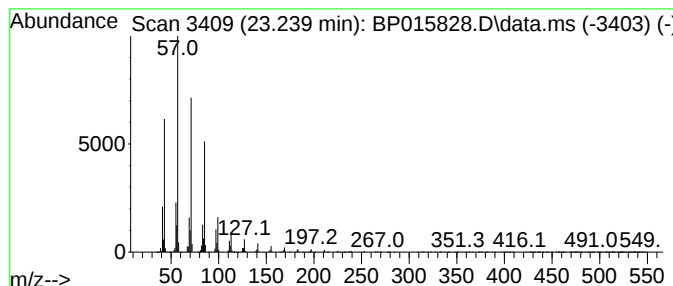
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Quant Title : SVOA CALIBRATION

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

Peak Number 18 (DEL) Alkane: Straight-Chai... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.239	35.75 ng/ul	5061070	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	97
2		Heptadecane, 9-octyl-	352	C25H52	007225-64-1	96
3		Nonadecane	268	C19H40	000629-92-5	93
4		Tetracosane	338	C24H50	000646-31-1	93
5		Eicosane	282	C20H42	000112-95-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFX5

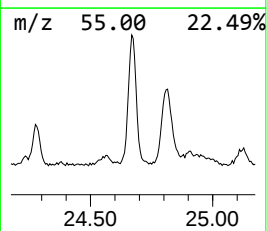
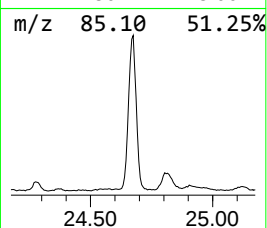
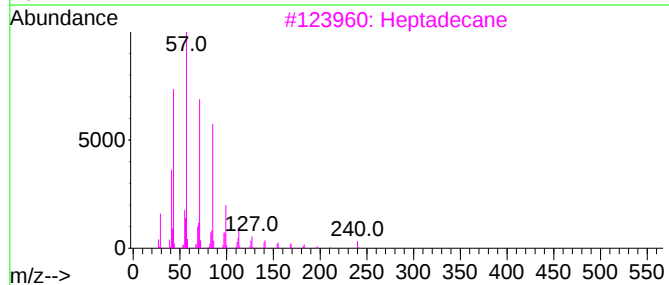
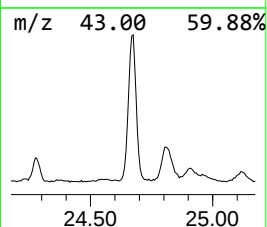
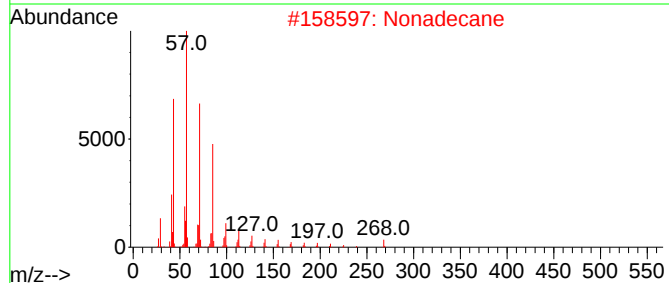
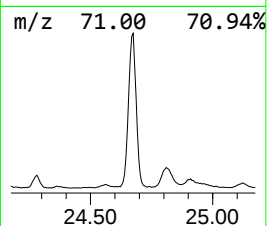
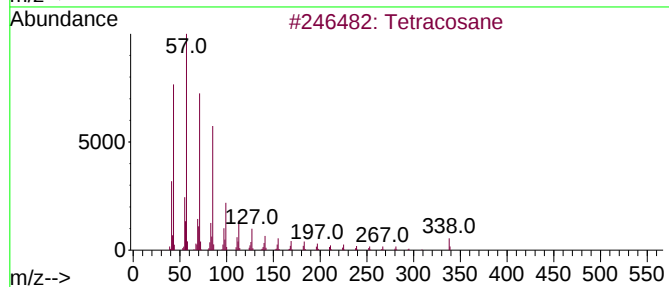
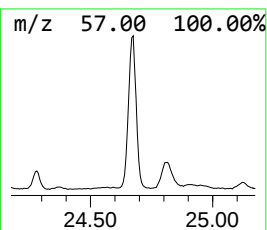
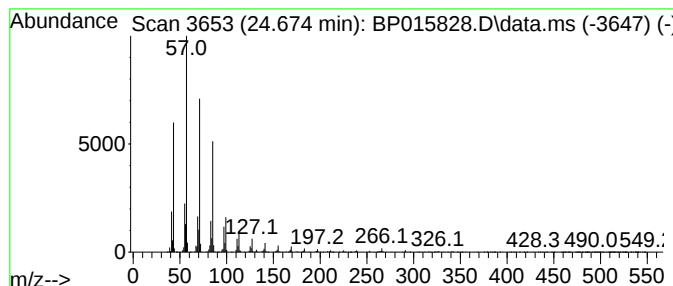
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Quant Title : SVOA CALIBRATION

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

Peak Number 19 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.674	26.98 ng/ul	3819740	Perylene-d12	24.098

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Tetracosane	338	C24H50	000646-31-1	96
2		Nonadecane	268	C19H40	000629-92-5	94
3		Heptadecane	240	C17H36	000629-78-7	93
4		Hexacosane	366	C26H54	000630-01-3	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
Data File : BP015828.D
Acq On : 30 Jun 2023 01:25
Operator : MA/JU
Sample : 03161-19
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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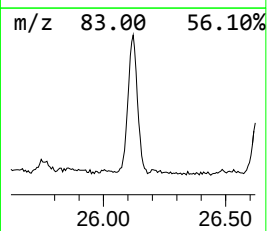
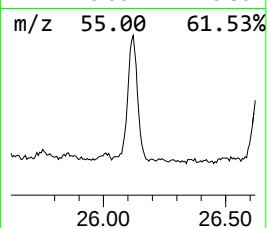
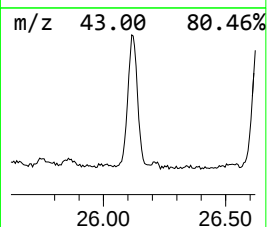
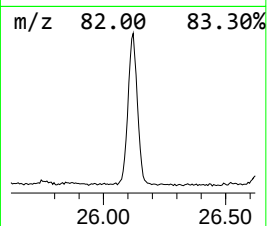
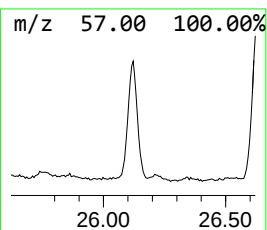
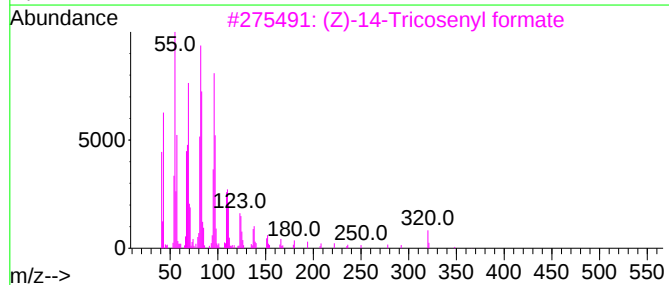
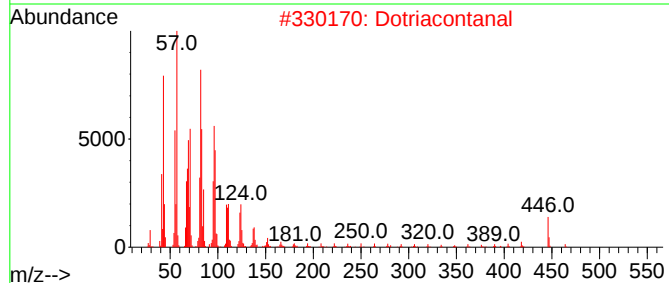
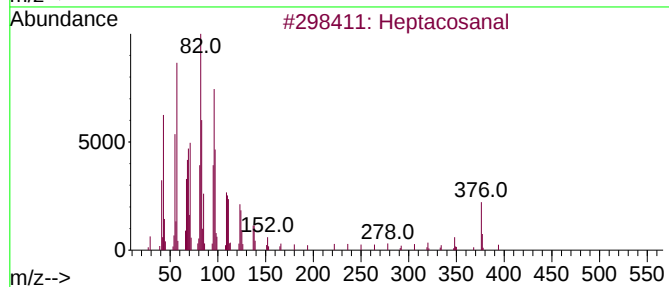
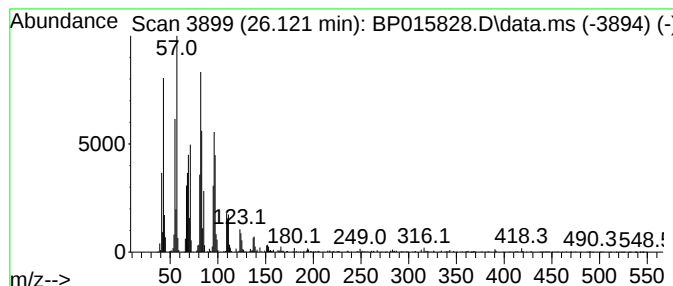
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Quant Title : SVOA CALIBRATION

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

Peak Number 20 Heptacosanal Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.121	18.07 ng/ul	2558150	Perylene-d12	24.098

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptacosanal	394	C27H54O	072934-03-3	97
2			Dotriacontanal	464	C32H64O	057878-00-9	91
3			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	83
4			Oxirane, heptadecyl-	282	C19H38O	067860-04-2	81
5			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	81



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062923\
 Data File : BP015828.D
 Acq On : 30 Jun 2023 01:25
 Operator : MA/JU
 Sample : 03161-19
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062623.MA.M
 Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
2H-Pyran, tetra...	3.928	3.4	ng/ul	246559	1	8.057	1460180	20.0
Benzophenone	16.069	3.9	ng/ul	564134	3	14.704	2922140	20.0
n-Hexadecanoic ...	18.316	15.0	ng/ul	2454970	4	17.463	3265840	20.0
1,1'-Biphenyl, ...	18.504	2.6	ng/ul	428312	4	17.463	3265840	20.0
1,1'-Biphenyl, ...	19.310	5.3	ng/ul	868421	4	17.463	3265840	20.0
Octadecanoic acid	19.539	7.6	ng/ul	1052770	5	21.551	2757880	20.0
1,1'-Biphenyl, ...	19.586	6.5	ng/ul	894540	5	21.551	2757880	20.0
1,1'-Biphenyl, ...	19.916	2.8	ng/ul	383215	5	21.551	2757880	20.0
1,1'-Biphenyl, ...	20.022	3.9	ng/ul	539389	5	21.551	2757880	20.0
unknown-01	20.269	5.0	ng/ul	687793	5	21.551	2757880	20.0
3,3',4,4',5-Pe...	20.327	3.5	ng/ul	488132	5	21.551	2757880	20.0
2,3,3',4',5,5'-...	20.539	2.6	ng/ul	353468	5	21.551	2757880	20.0
1,1'-Biphenyl, ...	20.598	2.9	ng/ul	393065	5	21.551	2757880	20.0
2,2',3,4,4',6'-...	20.851	2.7	ng/ul	367112	5	21.551	2757880	20.0
1-Heneicosyl fo...	21.221	17.4	ng/ul	2402420	5	21.551	2757880	20.0
(DEL) Alkane: S...	22.133	5.1	ng/ul	702887	5	21.551	2757880	20.0
1-Heneicosanol	22.174	15.9	ng/ul	2187660	5	21.551	2757880	20.0
(DEL) Alkane: S...	23.239	35.8	ng/ul	5061070	6	24.098	2831680	20.0
(DEL) Alkane: S...	24.674	27.0	ng/ul	3819740	6	24.098	2831680	20.0
Heptacosanal	26.121	18.1	ng/ul	2558150	6	24.098	2831680	20.0