

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
 Data File : BP015646.D
 Acq On : 22 Jun 2023 14:32
 Operator : MA/JU
 Sample : 03083-08
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DCFK6

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: BP015646.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.399	32	36	42	rVB	18344	25725	0.37%	0.071%
2	3.852	110	113	125	rBV	103885	170399	2.44%	0.473%
3	3.952	126	130	135	rVB	29532	39830	0.57%	0.111%
4	4.070	146	150	155	rVB	17791	27580	0.40%	0.077%
5	4.170	164	167	173	rVB2	9759	14514	0.21%	0.040%
6	6.046	483	486	490	rVB6	4977	7113	0.10%	0.020%
7	6.299	523	529	534	rBV5	9739	21059	0.30%	0.058%
8	7.240	683	689	706	rBV	276657	548965	7.88%	1.525%
9	7.399	709	716	728	rVV	236999	413208	5.93%	1.148%
10	7.499	730	733	740	rVB9	5745	9241	0.13%	0.026%
11	7.605	745	751	761	rBV	327888	564034	8.09%	1.567%
12	7.675	761	763	768	rVB2	9065	11685	0.17%	0.032%
13	8.075	825	831	841	rVV	504028	860963	12.35%	2.391%
14	8.805	948	955	973	rBV	311111	635338	9.12%	1.765%
15	9.257	1025	1032	1044	rBV	317423	599835	8.61%	1.666%
16	9.416	1053	1059	1066	rVB10	6533	15739	0.23%	0.044%
17	9.628	1090	1095	1100	rVB7	6141	12229	0.18%	0.034%
18	9.987	1149	1156	1167	rBV	259538	492150	7.06%	1.367%
19	10.540	1242	1250	1270	rBV	297462	730747	10.48%	2.030%
20	10.904	1305	1312	1319	rBV	725143	1253604	17.99%	3.482%
21	11.063	1332	1339	1358	rBV	127126	360395	5.17%	1.001%
22	11.751	1449	1456	1459	rBV9	4578	10151	0.15%	0.028%
23	11.904	1476	1482	1490	rBV5	10216	18262	0.26%	0.051%
24	12.310	1544	1551	1560	rVB5	7389	16924	0.24%	0.047%
25	12.510	1579	1585	1592	rBV9	6133	14213	0.20%	0.039%
26	12.575	1592	1596	1601	rBV5	6338	12910	0.19%	0.036%
27	12.787	1626	1632	1637	rBV6	7805	13279	0.19%	0.037%
28	13.098	1679	1685	1689	rVB8	2855	7559	0.11%	0.021%
29	13.240	1704	1709	1713	rVV4	8413	14724	0.21%	0.041%
30	13.528	1752	1758	1763	rBV3	12837	21499	0.31%	0.060%
31	13.604	1763	1771	1776	rVB4	15810	30988	0.44%	0.086%
32	13.710	1783	1789	1794	rVB6	12523	20493	0.29%	0.057%
33	13.893	1816	1820	1829	rBV6	6582	19696	0.28%	0.055%
34	14.045	1841	1846	1850	rVV6	10796	18707	0.27%	0.052%
35	14.134	1852	1861	1867	rVV	828230	1196611	17.17%	3.323%
36	14.181	1867	1869	1874	rVV5	21912	30160	0.43%	0.084%

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37	14.245	1874	1880	1884	rVV7	9730	17930	0.26%	0.050%
38	14.310	1889	1891	1896	rVB3	10529	12514	0.18%	0.035%
39	14.422	1903	1910	1922	rBV	872559	1359528	19.51%	3.776%
40	14.598	1936	1940	1945	rVB3	13939	19674	0.28%	0.055%

41	14.728	1955	1962	1970	rBV	1099901	1633929	23.44%	4.538%
42	14.798	1970	1974	1981	rVB9	17164	34143	0.49%	0.095%
43	14.916	1989	1994	1999	rBV	133663	236934	3.40%	0.658%
44	14.969	1999	2003	2024	rVB2	96090	300076	4.31%	0.833%
45	15.140	2027	2032	2034	rBV6	10808	17023	0.24%	0.047%

46	15.334	2060	2065	2073	rBV6	8376	24113	0.35%	0.067%
47	15.398	2073	2076	2081	rVB7	7426	11436	0.16%	0.032%
48	15.504	2091	2094	2099	rBV7	8915	13952	0.20%	0.039%
49	15.551	2099	2102	2107	rVB3	19242	22780	0.33%	0.063%
50	15.722	2124	2131	2146	rBV	1099656	1660625	23.83%	4.612%

51	15.863	2150	2155	2162	rBV	143187	286985	4.12%	0.797%
52	16.087	2187	2193	2204	rVB	410312	615384	8.83%	1.709%
53	16.222	2213	2216	2219	rVB5	7155	7743	0.11%	0.022%
54	16.304	2228	2230	2235	rVB5	7900	10981	0.16%	0.030%
55	16.434	2246	2252	2261	rBV5	34421	88833	1.27%	0.247%

56	16.587	2273	2278	2283	rBV7	13922	26106	0.37%	0.073%
57	16.651	2285	2289	2295	rVB	57835	84057	1.21%	0.233%
58	16.881	2324	2328	2331	rBV6	7285	12797	0.18%	0.036%
59	17.151	2371	2374	2377	rBV5	9248	13278	0.19%	0.037%
60	17.216	2381	2385	2390	rVB3	26601	34089	0.49%	0.095%

61	17.369	2406	2411	2417	rBV	51482	92157	1.32%	0.256%
62	17.434	2419	2422	2426	rVV5	26661	42097	0.60%	0.117%
63	17.486	2426	2431	2436	rVV	1177656	1706124	24.48%	4.739%
64	17.534	2436	2439	2443	rVV	90485	131196	1.88%	0.364%
65	17.586	2443	2448	2457	rVB2	922015	1401541	20.11%	3.893%

66	17.951	2507	2510	2515	rVB	30366	40093	0.58%	0.111%
67	18.239	2556	2559	2562	rVB3	21944	22408	0.32%	0.062%
68	18.292	2562	2568	2573	rBV4	70563	131412	1.89%	0.365%
69	18.345	2573	2577	2584	rBV	429511	665729	9.55%	1.849%
70	18.622	2621	2624	2634	rBV6	33904	90466	1.30%	0.251%

71	19.228	2725	2727	2733	rVB3	63483	83481	1.20%	0.232%
72	19.328	2741	2744	2748	rVB5	33990	38309	0.55%	0.106%
73	19.392	2751	2755	2759	rVV2	97071	119000	1.71%	0.331%
74	19.439	2760	2763	2765	rBV2	42580	53842	0.77%	0.150%
75	19.492	2768	2772	2779	rVB	255912	370156	5.31%	1.028%

76	19.580	2783	2787	2791	rBV2	134911	211056	3.03%	0.586%
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77	19.622	2791	2794	2801	rVB5	84494	126601	1.82%	0.352%
78	19.780	2817	2821	2822	rBV4	124934	146716	2.10%	0.407%
79	19.816	2822	2827	2831	rVV	1086904	1488205	21.35%	4.133%
80	20.127	2877	2880	2886	rVB6	114549	162480	2.33%	0.451%
81	20.751	2983	2986	2989	rVB	212662	209022	3.00%	0.581%
82	21.251	3067	3071	3080	rBV2	276822	446715	6.41%	1.241%
83	21.580	3121	3127	3131	rBV	978323	1494047	21.44%	4.150%
84	23.292	3414	3418	3423	rVB	396894	614119	8.81%	1.706%
85	23.963	3526	3532	3539	rBV2	859385	1655493	23.75%	4.598%
86	24.133	3555	3561	3570	rVB	669253	1396210	20.03%	3.878%
87	24.745	3660	3665	3676	rVB	585527	1291082	18.52%	3.586%
88	28.874	4360	4367	4384	rVB3	1866383	6969897	100.00%	19.358%

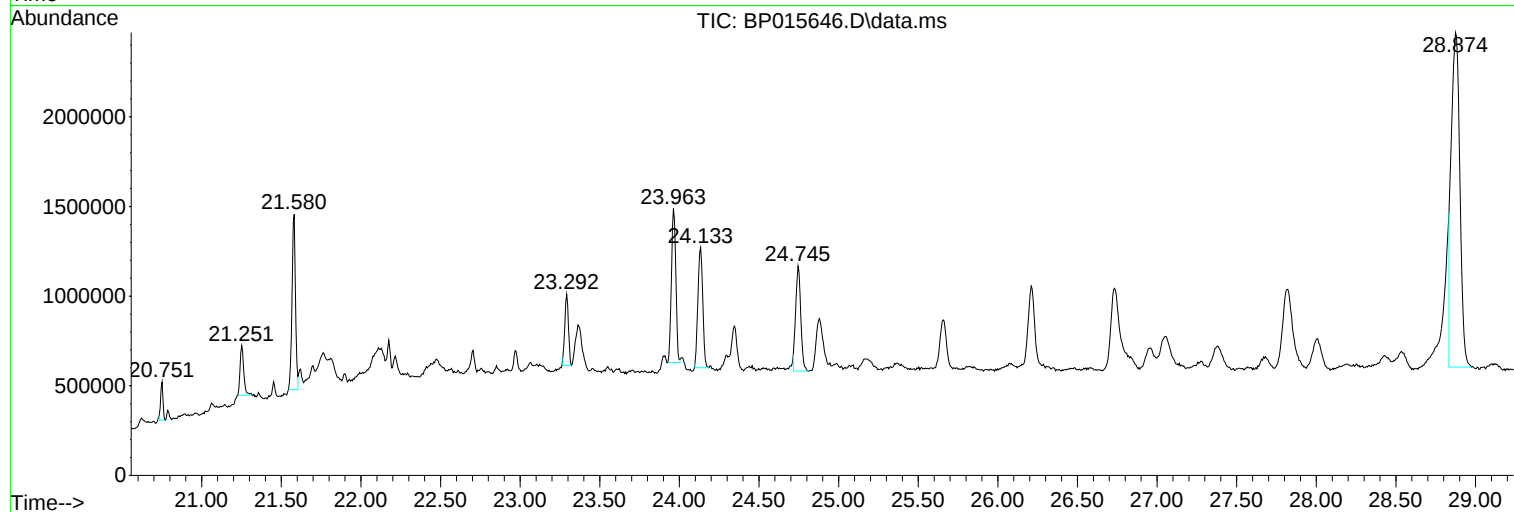
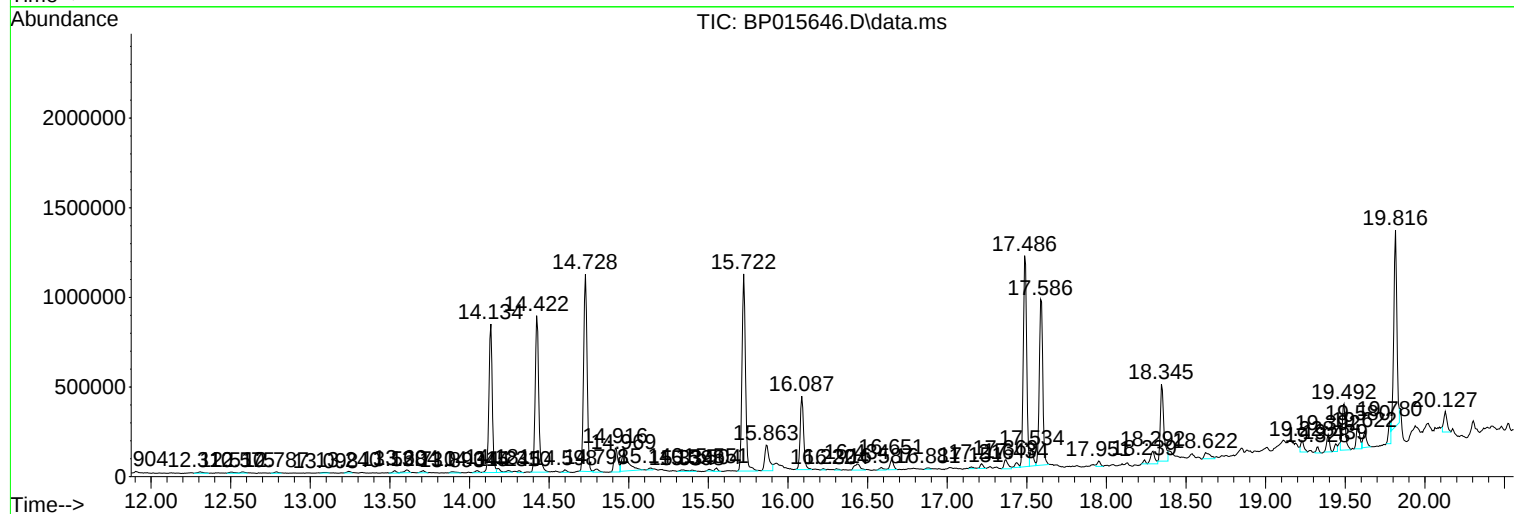
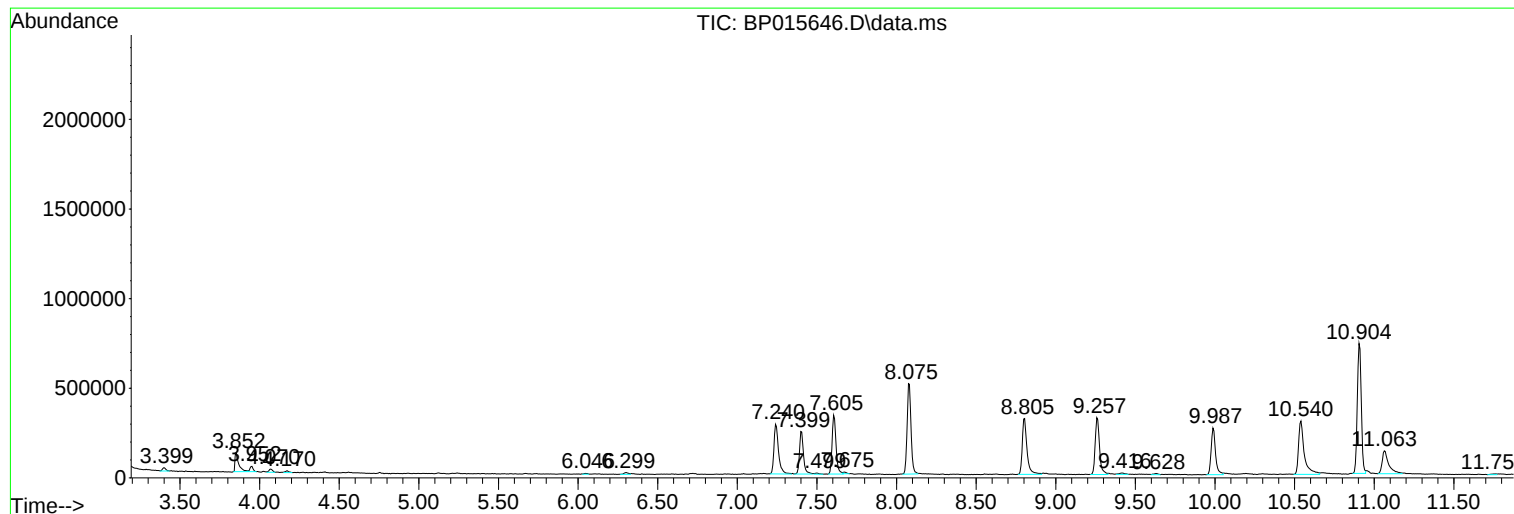
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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



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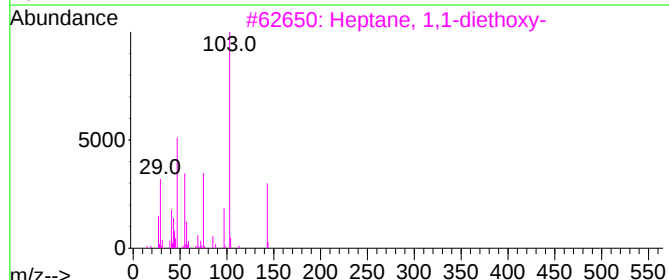
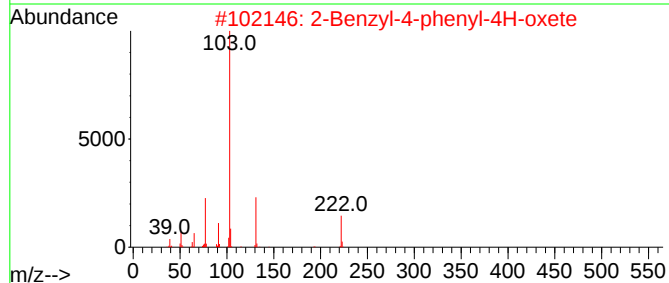
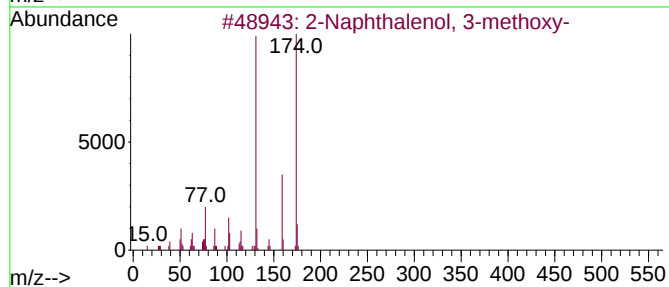
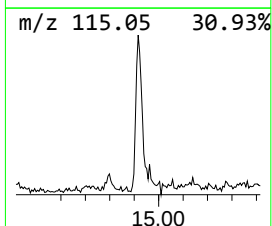
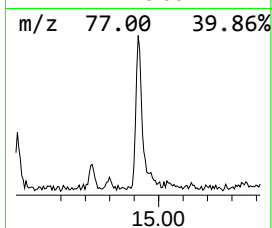
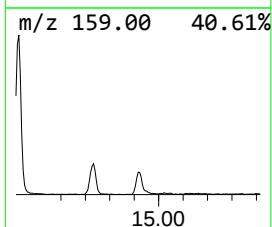
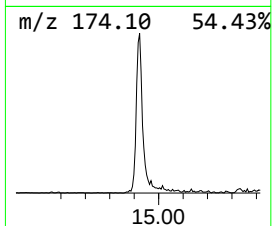
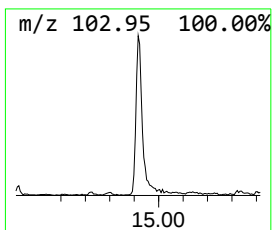
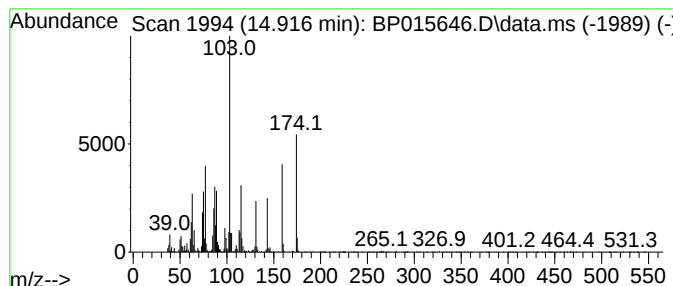
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 unknown-01 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.916	2.90 ng/ul	236934	Acenaphthene-d10	14.728

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenol, 3-methoxy-		174	C11H10O2		018515-11-2	18
2	2-Benzyl-4-phenyl-4H-oxete		222	C16H14O		1000444-12-8	14
3	Heptane, 1,1-diethoxy-		188	C11H24O2		000688-82-4	14
4	3H-Pyrazol-3-one, 2,4-dihydro-2-...		174	C10H10N2O		041927-50-8	11
5	2H-1-Benzopyran-2-one 3,4-dimethyl-		174	C11H10O2		004281-39-4	11



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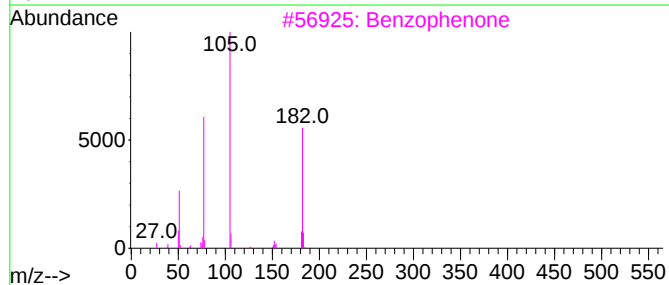
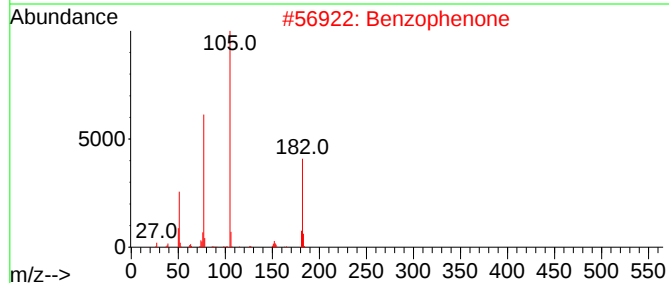
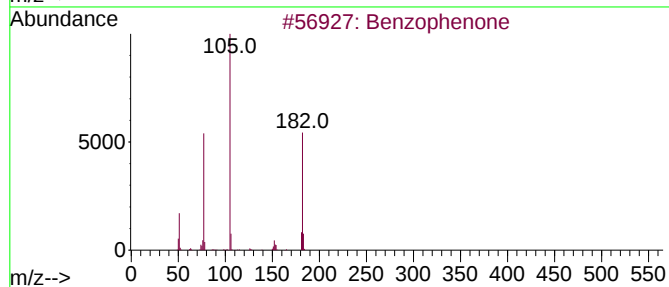
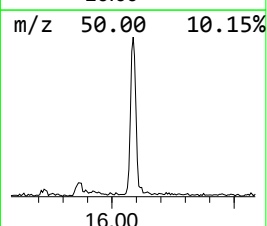
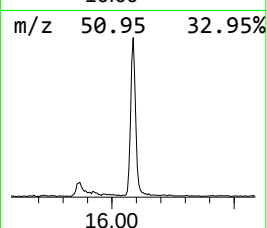
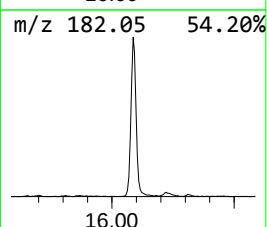
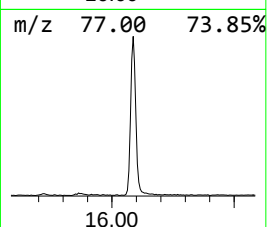
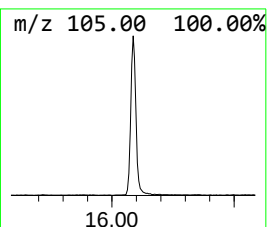
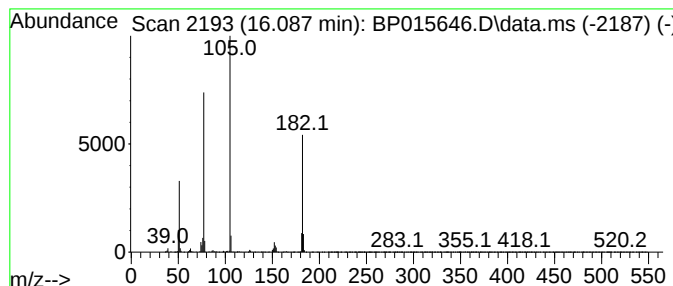
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzophenone Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.087	7.53 ng/ul	615384	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	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	95
5			Benzophenone	182	C13H10O	000119-61-9	91



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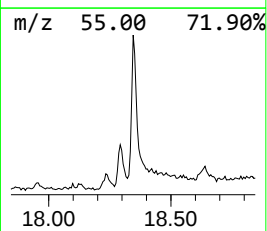
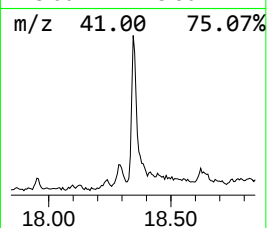
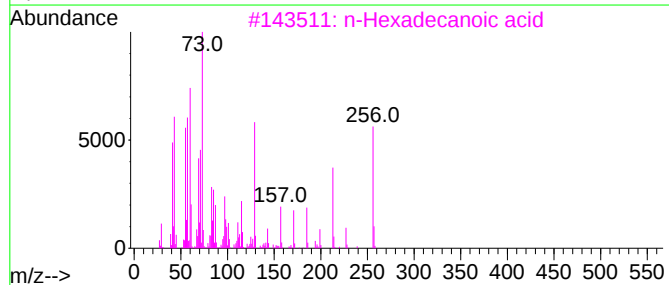
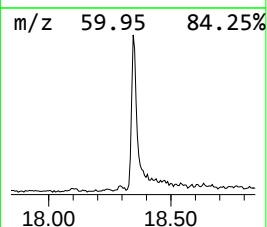
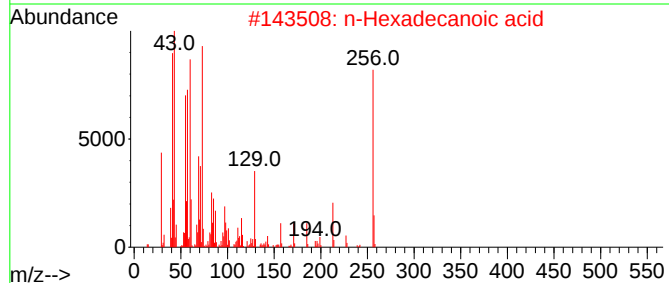
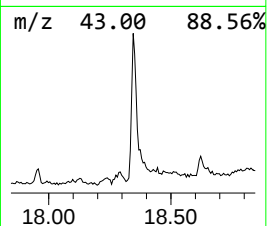
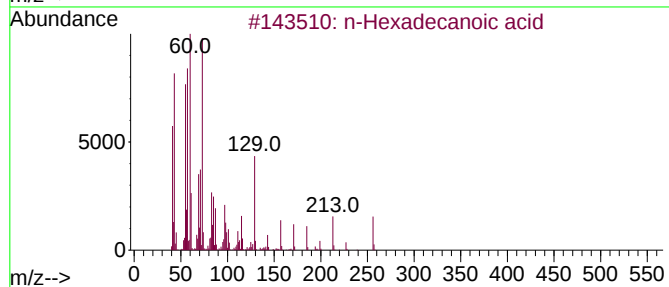
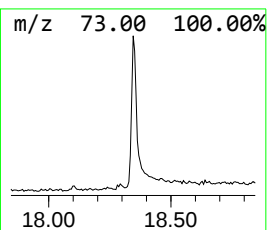
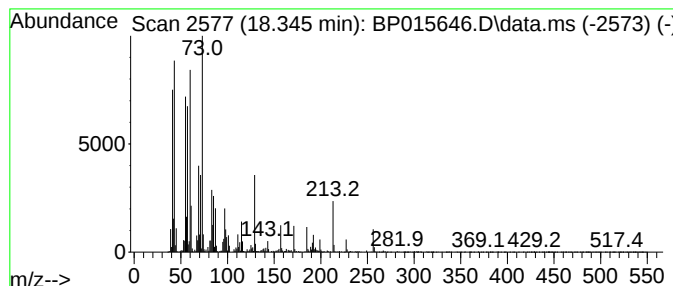
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.345	7.80 ng/ul	665729	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	97
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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ClientSampleId :
DCFK6

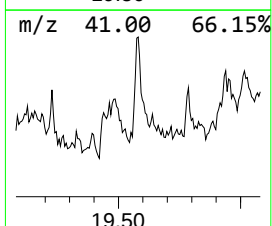
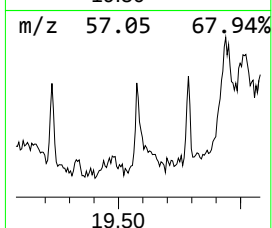
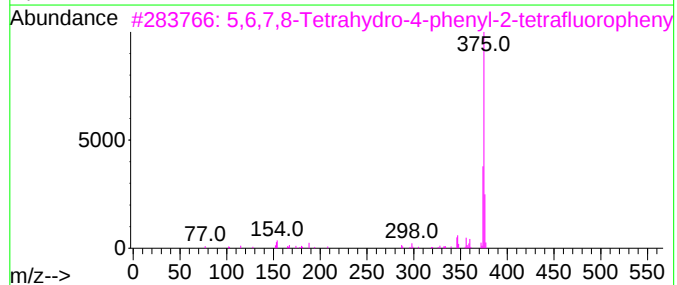
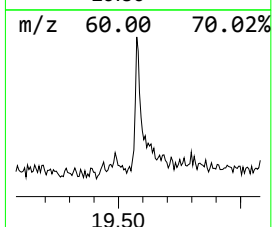
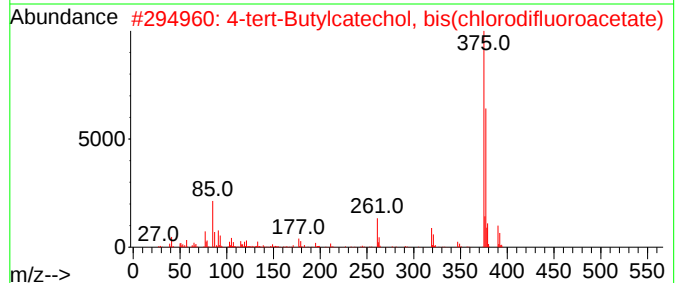
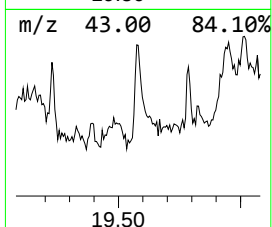
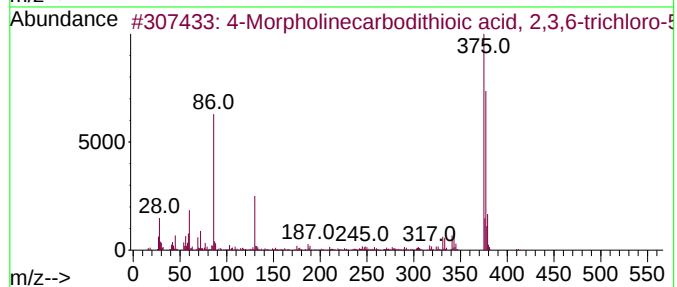
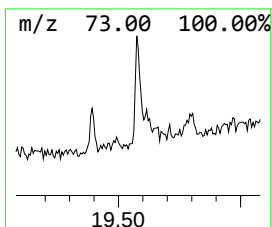
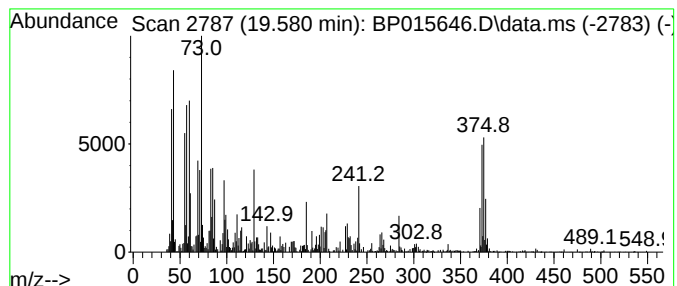
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-02 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.581	2.83 ng/ul	211056	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	4-Morpholinecarbodithioic acid, ...	410	C11H8Cl3F3N2OS2	1000305-31-7	10		
2	4-tert-Butylcatechol, bis(chloro...	390	C14H12Cl2F4O4	1000447-47-8	10		
3	5,6,7,8-Tetrahydro-4-phenyl-2-te...	375	C21H14F5N	140484-19-1	10		
4	Indole-3-carboxylic acid, 5-acet...	415	C20H18BrNO4	040945-73-1	9		
5	1-Monopalmitin, 2TMS derivative	474	C25H54O4Si2	001188-74-5	9		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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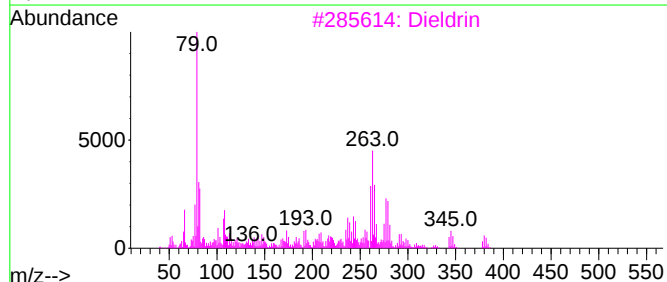
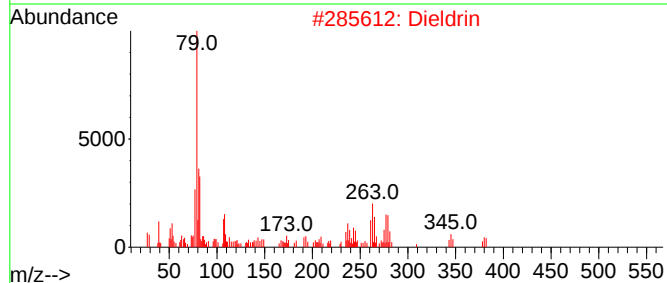
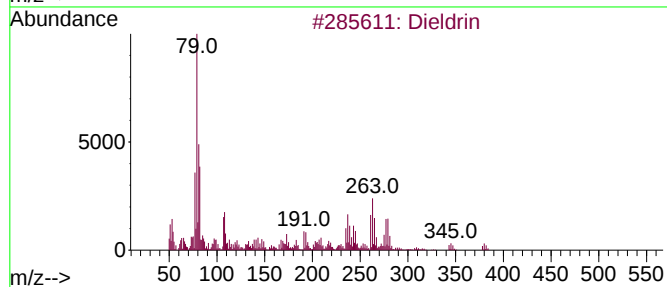
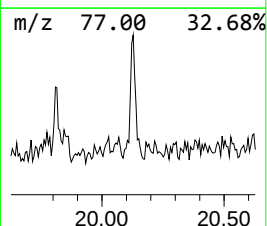
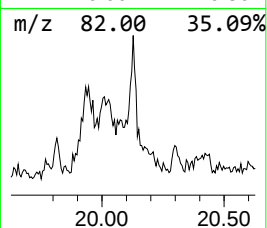
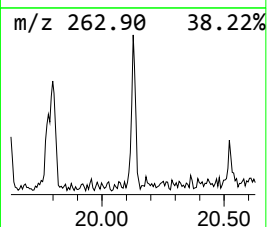
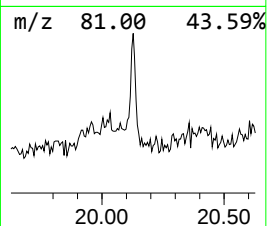
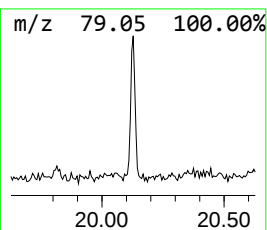
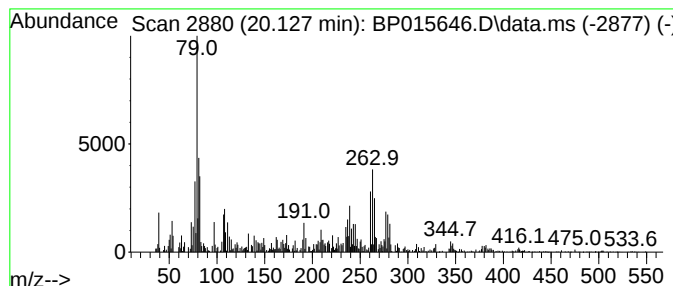
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Dieldrin Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.128	2.18 ng/ul	162480	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dieldrin	378	C12H8Cl6O	000060-57-1	98
2			Dieldrin	378	C12H8Cl6O	000060-57-1	95
3			Dieldrin	378	C12H8Cl6O	000060-57-1	93
4			Dieldrin	378	C12H8Cl6O	000060-57-1	86
5			Dieldrin	378	C12H8Cl6O	000060-57-1	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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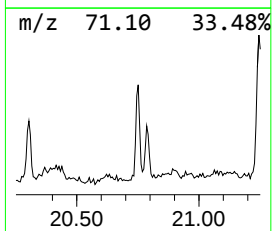
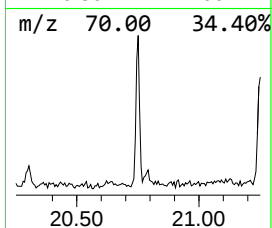
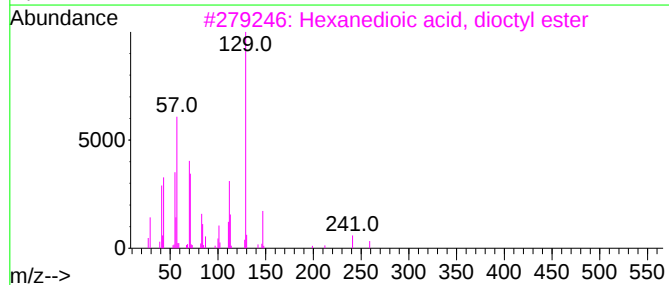
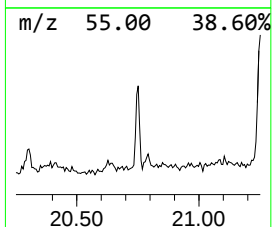
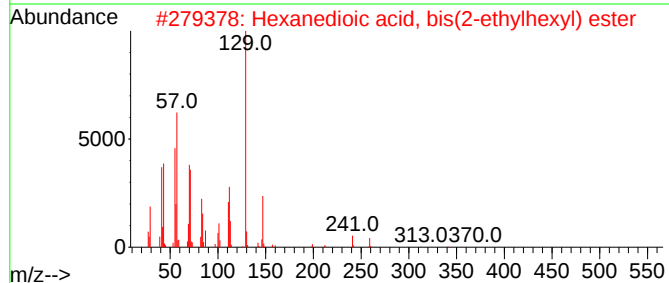
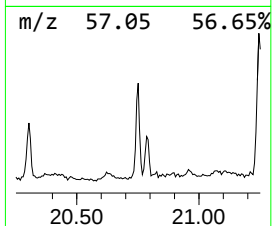
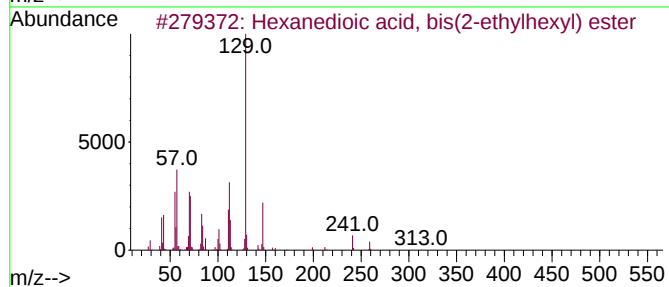
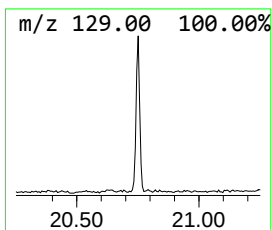
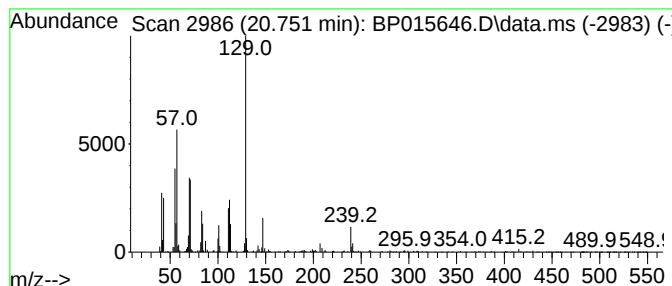
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Hexanedioic acid, bis(2-eth... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.751	2.80 ng/ul	209022	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	90
2			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	83
3			Hexanedioic acid, dioctyl ester	370	C22H42O4	000123-79-5	81
4			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	81
5			Hexanedioic acid, bis(2-ethylhex...	370	C22H42O4	000103-23-1	74



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK6

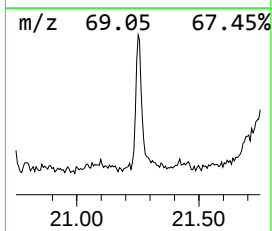
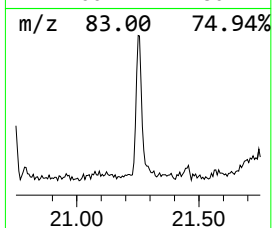
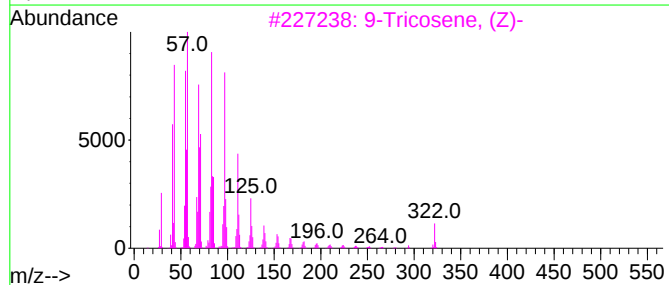
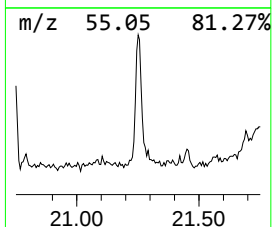
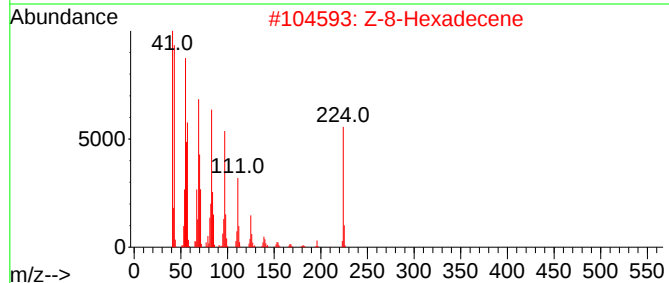
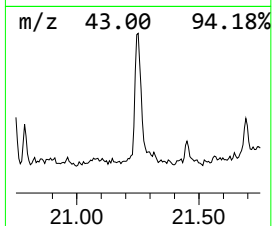
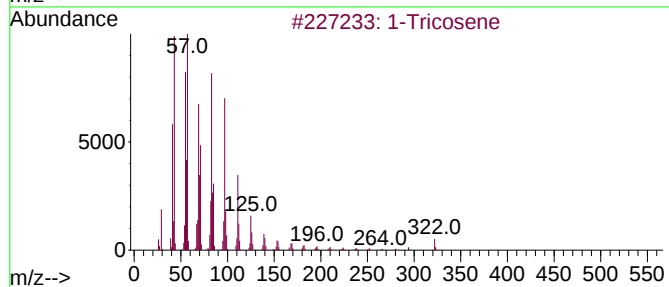
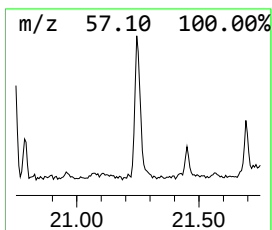
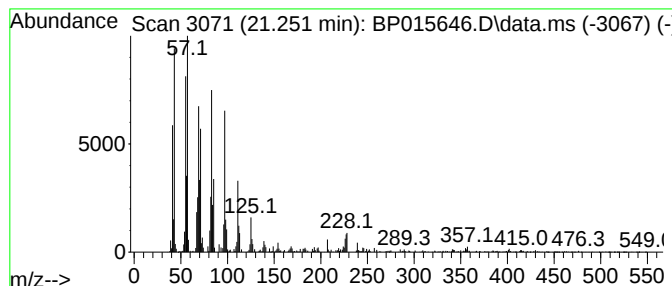
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.251	5.98 ng/ul	446715	Chrysene-d12	21.580

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Tricosene	322	C23H46	018835-32-0	96
2			Z-8-Hexadecene	224	C16H32	1000130-87-5	96
3			9-Tricosene, (Z)-	322	C23H46	027519-02-4	95
4			E-14-Hexadecenal	238	C16H30O	330207-53-9	94
5			Pentadecafluorooctanoic acid, oc...	666	C26H37F15O2	1000406-04-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

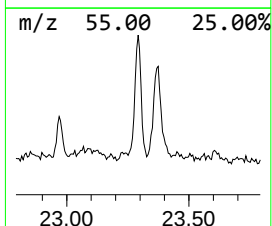
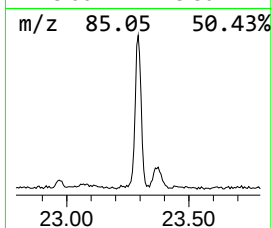
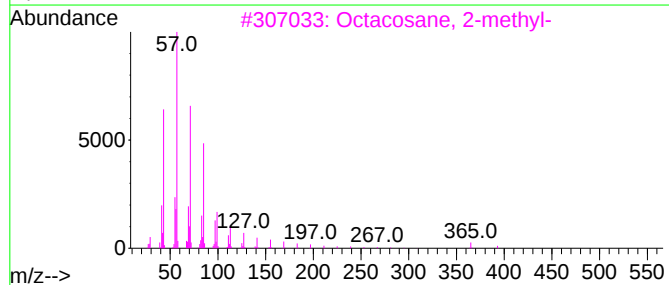
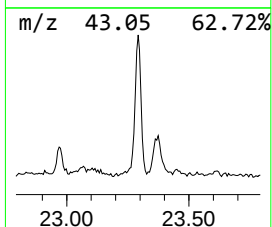
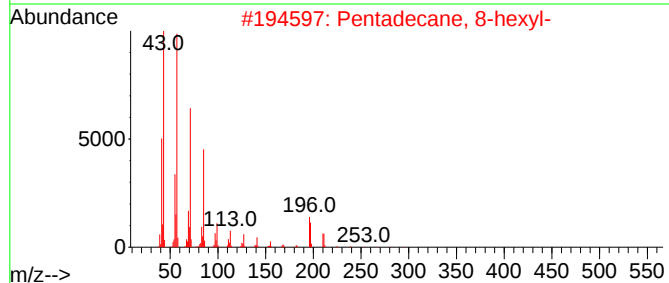
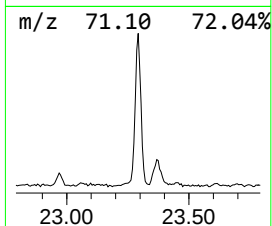
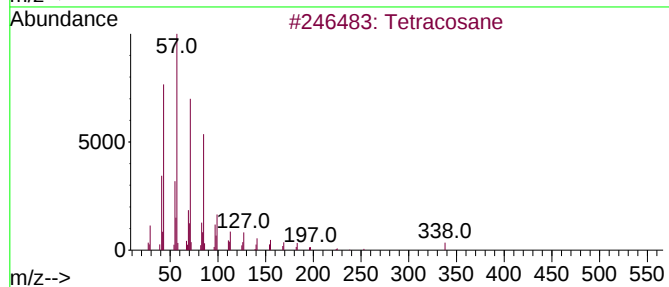
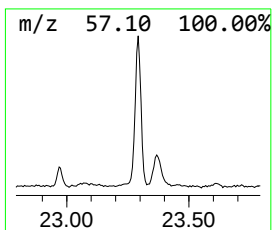
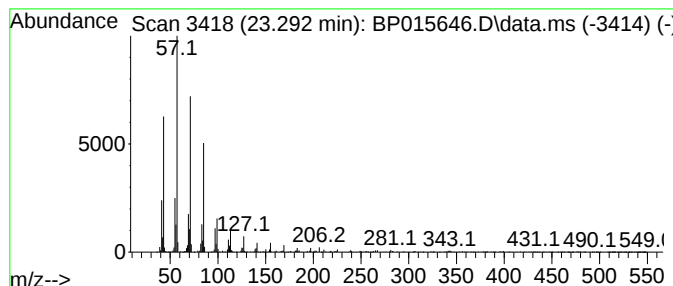
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
23.292	8.80 ng/ul	614119	Perylene-d12	24.133	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Tetracosane	338	C24H50	000646-31-1	93
2	Pentadecane, 8-hexyl-	296	C21H44	013475-75-7	93
3	Octacosane, 2-methyl-	408	C29H60	001560-98-1	90
4	Heptadecane, 9-octyl-	352	C25H52	007225-64-1	90
5	Octacosane	394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

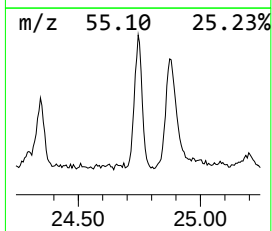
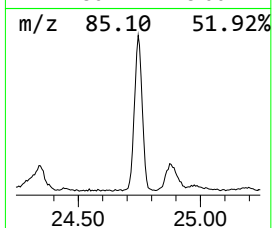
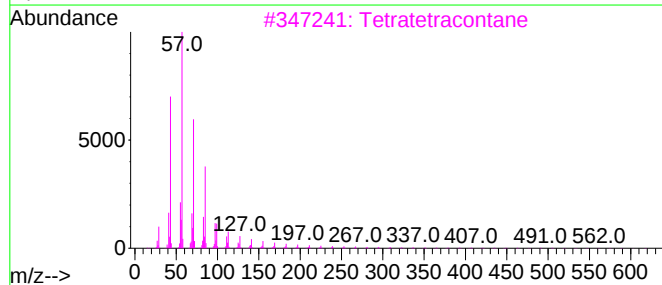
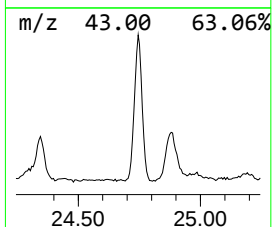
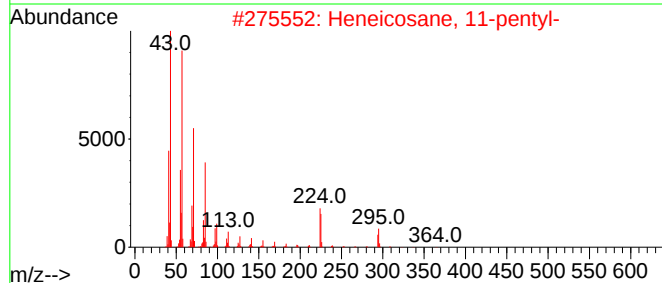
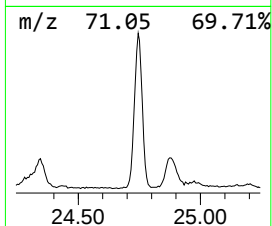
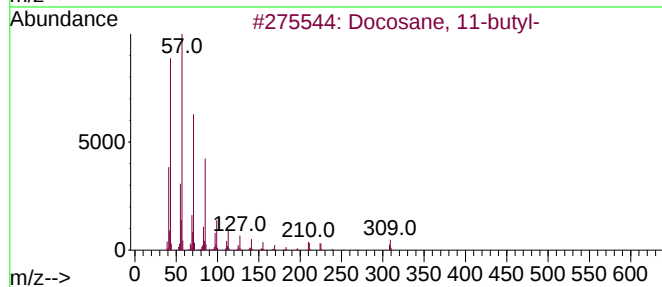
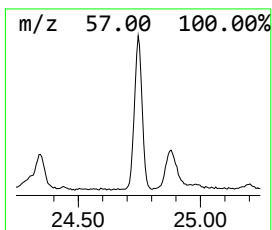
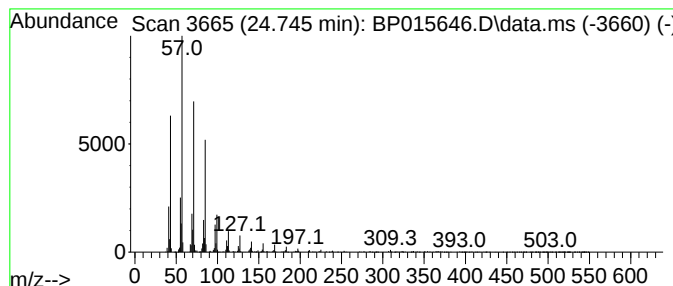
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.745	18.49 ng/ul	1291080	Perylene-d12	24.133	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Docosane, 11-butyl-	366	C26H54	013475-76-8	95
2	Heneicosane, 11-pentyl-	366	C26H54	014739-72-1	93
3	Tetratetracontane	619	C44H90	007098-22-8	91
4	Heneicosane	296	C21H44	000629-94-7	91
5	Octacosane, 1-iodo-	520	C28H57I	1000406-32-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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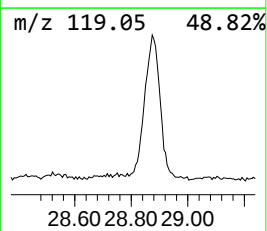
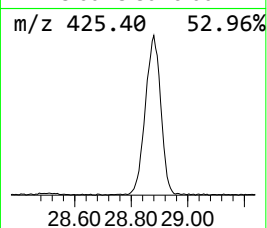
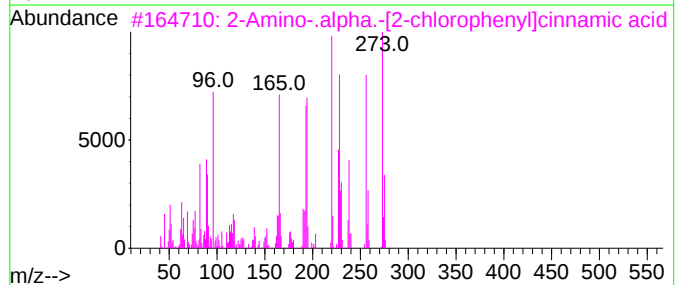
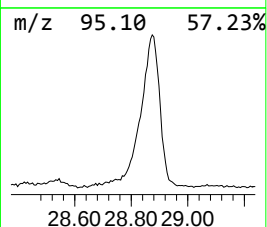
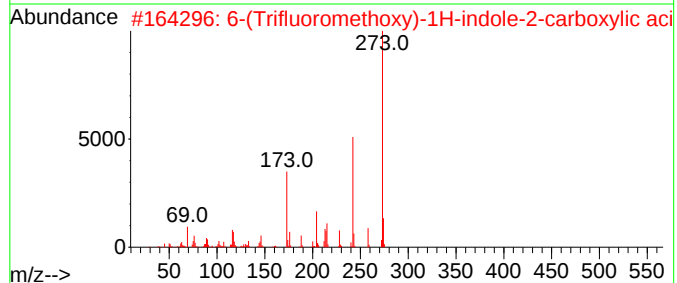
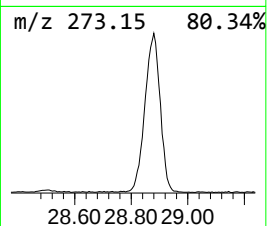
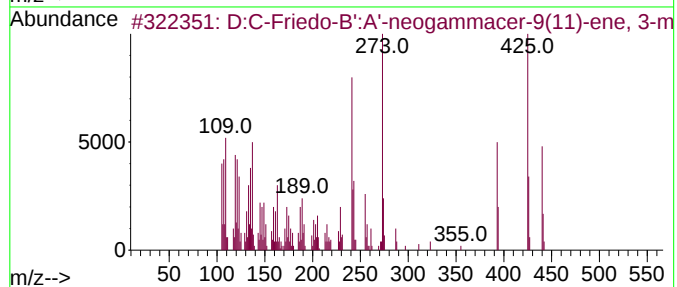
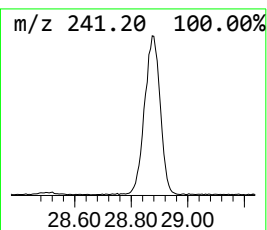
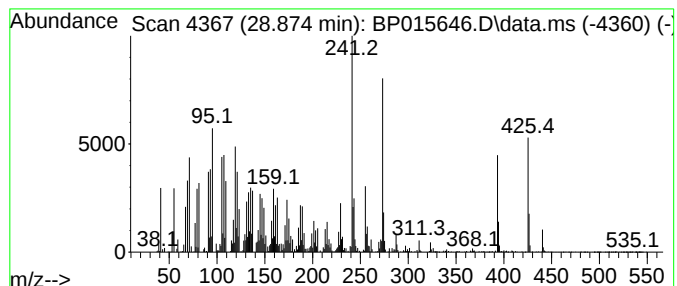
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-03 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.874	99.84 ng/ul	6969900	Perylene-d12	24.133

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	46
2			6-(Trifluoromethoxy)-1H-indole-2...	273	C12H10F3NO3	1000499-75-7	25
3			2-Amino-.alpha.-[2-chlorophenyl]...	273	C15H12ClNO2	1000254-08-5	25
4			1,3-Dibromo-5,7-dimethyladamantane	320	C12H18Br2	1000411-41-6	22
5			3-Bromo-4-hydroxy-5-methoxypheny...	241	C9H8BrNO2	081038-44-0	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062223\
Data File : BP015646.D
Acq On : 22 Jun 2023 14:32
Operator : MA/JU
Sample : 03083-08
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DCFK6

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
unknown-01	14.916	2.9	ng/ul	236934	3	14.728	1633930	20.0
Benzophenone	16.087	7.5	ng/ul	615384	3	14.728	1633930	20.0
n-Hexadecanoic ...	18.345	7.8	ng/ul	665729	4	17.486	1706120	20.0
unknown-02	19.581	2.8	ng/ul	211056	5	21.580	1494050	20.0
Dieldrin	20.128	2.2	ng/ul	162480	5	21.580	1494050	20.0
Hexanedioic aci...	20.751	2.8	ng/ul	209022	5	21.580	1494050	20.0
(DEL) Alkane: S...	21.251	6.0	ng/ul	446715	5	21.580	1494050	20.0
(DEL) Alkane: S...	23.292	8.8	ng/ul	614119	6	24.133	1396210	20.0
(DEL) Alkane: S...	24.745	18.5	ng/ul	1291080	6	24.133	1396210	20.0
unknown-03	28.874	99.8	ng/ul	6969900	6	24.133	1396210	20.0