

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
 Data File : BP015403.D
 Acq On : 07 Jun 2023 17:34
 Operator : MA/JU
 Sample : 02950-06
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EW983

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: BP015403.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.411	35	38	46	rVB2	16589	25315	0.74%	0.065%
2	3.858	112	114	118	rVB4	4381	5016	0.15%	0.013%
3	3.917	122	124	125	rBV2	6639	4168	0.12%	0.011%
4	3.964	129	132	137	rVB3	15594	24116	0.71%	0.062%
5	4.087	148	153	159	rVB3	10352	18272	0.54%	0.047%
6	4.187	166	170	176	rBV	11882	19852	0.58%	0.051%
7	4.528	226	228	233	rBV6	2684	4037	0.12%	0.010%
8	4.581	233	237	242	rVV7	4799	9183	0.27%	0.024%
9	4.640	243	247	252	rVB2	7783	12422	0.36%	0.032%
10	5.158	328	335	340	rBV4	8540	19608	0.58%	0.051%
11	5.505	390	394	396	rVB5	2988	3925	0.12%	0.010%
12	5.628	411	415	418	rVB5	2867	3687	0.11%	0.010%
13	6.052	482	487	493	rVB6	6452	13342	0.39%	0.034%
14	7.069	656	660	662	rBV4	3774	4198	0.12%	0.011%
15	7.199	674	682	684	rBV5	5678	10326	0.30%	0.027%
16	7.252	684	691	705	rVV	272378	498878	14.65%	1.287%
17	7.422	713	720	732	rVB	241426	387148	11.37%	0.999%
18	7.622	748	754	762	rBV	286435	476714	14.00%	1.230%
19	7.705	764	768	779	rVB	21337	48256	1.42%	0.125%
20	7.916	800	804	808	rVB7	4340	6836	0.20%	0.018%
21	8.016	816	821	826	rVB9	5689	10851	0.32%	0.028%
22	8.099	826	835	844	rBV	548524	870115	25.56%	2.245%
23	8.269	860	864	869	rBV8	3530	7090	0.21%	0.018%
24	8.640	921	927	933	rBV8	3399	7635	0.22%	0.020%
25	8.816	950	957	972	rBV	291710	538732	15.82%	1.390%
26	8.934	972	977	993	rVV	182051	322163	9.46%	0.831%
27	9.105	1002	1006	1011	rVV8	3375	5792	0.17%	0.015%
28	9.275	1028	1035	1045	rBV	302364	519077	15.25%	1.339%
29	9.446	1058	1064	1069	rBV10	4858	13528	0.40%	0.035%
30	9.569	1082	1085	1092	rVB9	3614	7008	0.21%	0.018%
31	9.663	1092	1101	1106	rBV	19218	37503	1.10%	0.097%
32	10.005	1152	1159	1169	rVV	259715	466141	13.69%	1.203%
33	10.093	1171	1174	1182	rVB8	9909	19680	0.58%	0.051%
34	10.228	1190	1197	1198	rBV7	4015	6680	0.20%	0.017%
35	10.375	1215	1222	1227	rVB3	15752	31894	0.94%	0.082%
36	10.493	1236	1242	1243	rBV6	7278	12317	0.36%	0.032%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
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37	10.552	1245	1252	1264	rVV	335074	641932	18.85%	1.656%
38	10.834	1296	1300	1301	rBV4	4074	4007	0.12%	0.010%
39	10.881	1303	1308	1309	rVV5	6146	7711	0.23%	0.020%
40	10.928	1309	1316	1325	rVV	806753	1356215	39.83%	3.500%
41	11.075	1334	1341	1356	rBV	231739	522340	15.34%	1.348%
42	11.405	1392	1397	1403	rVB4	25917	42807	1.26%	0.110%
43	11.469	1404	1408	1409	rBV3	6330	7197	0.21%	0.019%
44	11.481	1409	1410	1415	rVB5	8748	10727	0.32%	0.028%
45	11.581	1425	1427	1431	rBV4	6361	7963	0.23%	0.021%
46	11.681	1439	1444	1445	rBV5	4762	7252	0.21%	0.019%
47	11.752	1449	1456	1460	rBV8	17535	42728	1.25%	0.110%
48	11.828	1466	1469	1470	rBV3	7497	7156	0.21%	0.018%
49	11.893	1477	1480	1484	rBV6	5220	7147	0.21%	0.018%
50	11.969	1489	1493	1497	rVV3	26411	41830	1.23%	0.108%
51	12.010	1497	1500	1505	rVB7	12295	23262	0.68%	0.060%
52	12.110	1512	1517	1522	rBV7	12457	18950	0.56%	0.049%
53	12.163	1522	1526	1528	rBV3	11915	17378	0.51%	0.045%
54	12.193	1528	1531	1535	rVB2	23971	32915	0.97%	0.085%
55	12.516	1584	1586	1595	rVB10	19472	36339	1.07%	0.094%
56	12.640	1603	1607	1609	rBV5	11495	17703	0.52%	0.046%
57	12.728	1618	1622	1623	rBV4	14870	18419	0.54%	0.048%
58	12.952	1657	1660	1666	rVB7	16639	22595	0.66%	0.058%
59	13.087	1680	1683	1685	rBV4	10683	16007	0.47%	0.041%
60	13.181	1695	1699	1704	rVB6	15497	25255	0.74%	0.065%
61	13.546	1756	1761	1764	rBV6	38927	66057	1.94%	0.170%
62	13.628	1770	1775	1782	rBV6	30161	66439	1.95%	0.171%
63	13.787	1800	1802	1807	rVB6	18852	26244	0.77%	0.068%
64	13.934	1825	1827	1831	rVB4	16864	19146	0.56%	0.049%
65	14.057	1846	1848	1852	rBV4	15618	19590	0.58%	0.051%
66	14.146	1856	1863	1870	rBV	989146	1383219	40.63%	3.569%
67	14.440	1907	1913	1929	rBV	992385	1593110	46.79%	4.111%
68	14.569	1931	1935	1939	rVB4	31948	50625	1.49%	0.131%
69	14.746	1959	1965	1972	rVB	1309130	1970546	57.88%	5.085%
70	14.951	1995	2000	2010	rBV	163400	364912	10.72%	0.942%
71	15.734	2128	2133	2139	rBV	1331628	1963092	57.66%	5.066%
72	15.863	2150	2155	2164	rVB	298986	486996	14.30%	1.257%
73	16.098	2190	2195	2200	rVB	476619	630603	18.52%	1.627%
74	16.234	2215	2218	2221	rBV3	35585	49319	1.45%	0.127%
75	16.663	2288	2291	2296	rVB3	81733	106910	3.14%	0.276%
76	17.316	2400	2402	2408	rVB3	53416	68530	2.01%	0.177%

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77	17.498	2427	2433	2438	rBV	1787558	2530158	74.32%	6.529%
78	17.598	2445	2450	2458	rVB	1584074	2191636	64.37%	5.655%
79	17.687	2460	2465	2470	rBV3	102495	228708	6.72%	0.590%
80	17.734	2470	2473	2476	rVV2	77476	97384	2.86%	0.251%
81	17.781	2476	2481	2486	rVB4	77599	182075	5.35%	0.470%
82	18.204	2550	2553	2558	rBV	102999	132065	3.88%	0.341%
83	18.357	2575	2579	2592	rBV8	326170	707645	20.78%	1.826%
84	18.481	2598	2600	2605	rVB3	72074	96443	2.83%	0.249%
85	18.739	2640	2644	2649	rVB	243424	310456	9.12%	0.801%
86	18.828	2656	2659	2663	rBV2	190964	308945	9.07%	0.797%
87	18.922	2672	2675	2678	rVB2	169597	190997	5.61%	0.493%
88	19.239	2726	2729	2732	rBV2	107590	148220	4.35%	0.382%
89	19.492	2768	2772	2777	rBV2	500339	651521	19.14%	1.681%
90	19.551	2778	2782	2785	rVV2	213876	300117	8.81%	0.774%
91	19.581	2785	2787	2791	rVB3	161920	196001	5.76%	0.506%
92	19.822	2823	2828	2831	rBV	2147536	3078311	90.42%	7.944%
93	20.481	2937	2940	2944	rVB	360462	428233	12.58%	1.105%
94	20.616	2961	2963	2967	rBV	181151	202902	5.96%	0.524%
95	21.222	3063	3066	3069	rBV	434031	511871	15.03%	1.321%
96	21.257	3069	3072	3077	rVB2	396727	567858	16.68%	1.465%
97	21.580	3121	3127	3131	rBV	2201012	3404628	100.00%	8.786%
98	21.616	3131	3133	3143	rVB	616723	833439	24.48%	2.151%
99	23.969	3527	3533	3539	rBV2	1284232	2505648	73.60%	6.466%
100	24.133	3555	3561	3570	rVB	1291382	2674438	78.55%	6.901%

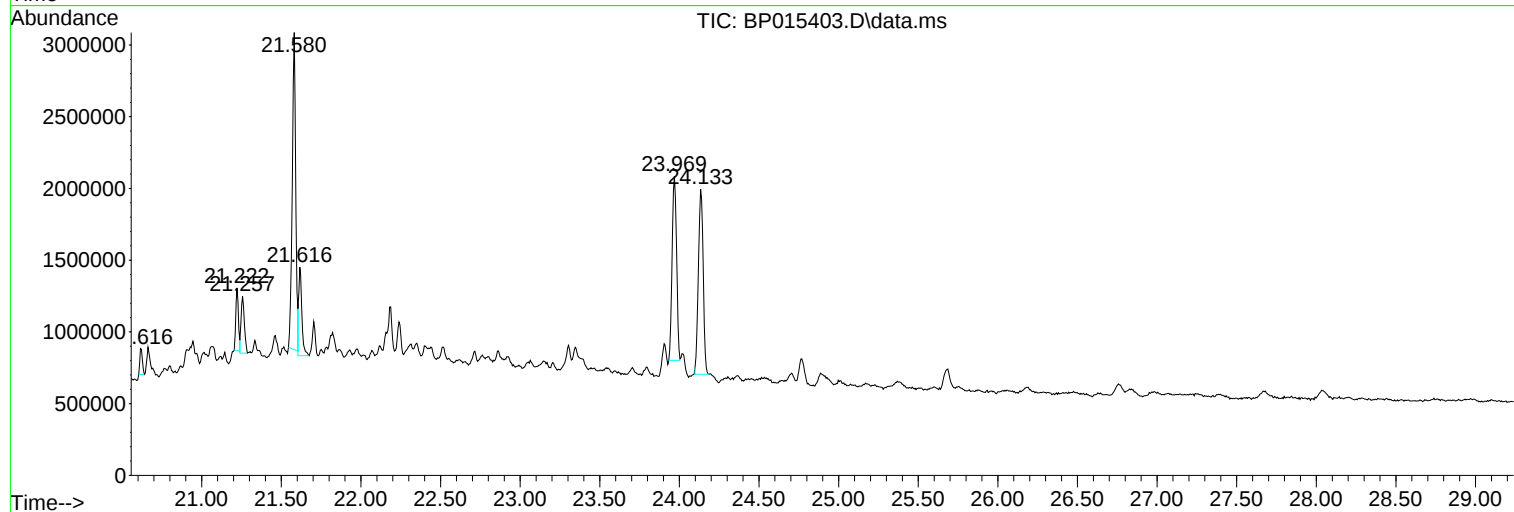
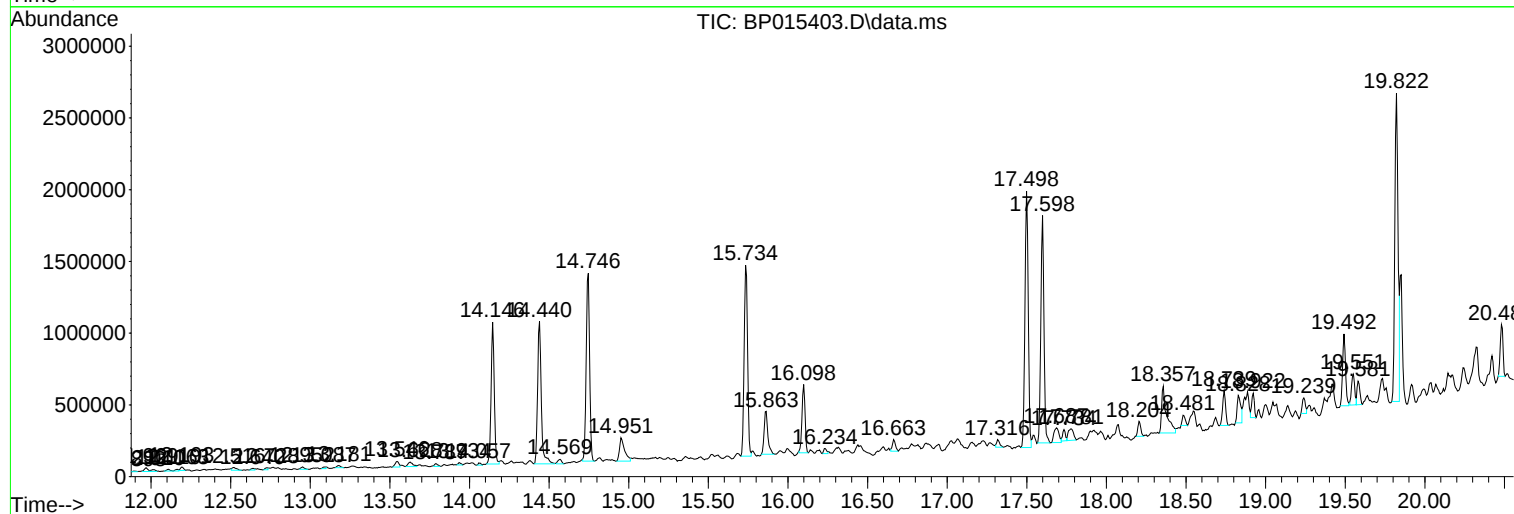
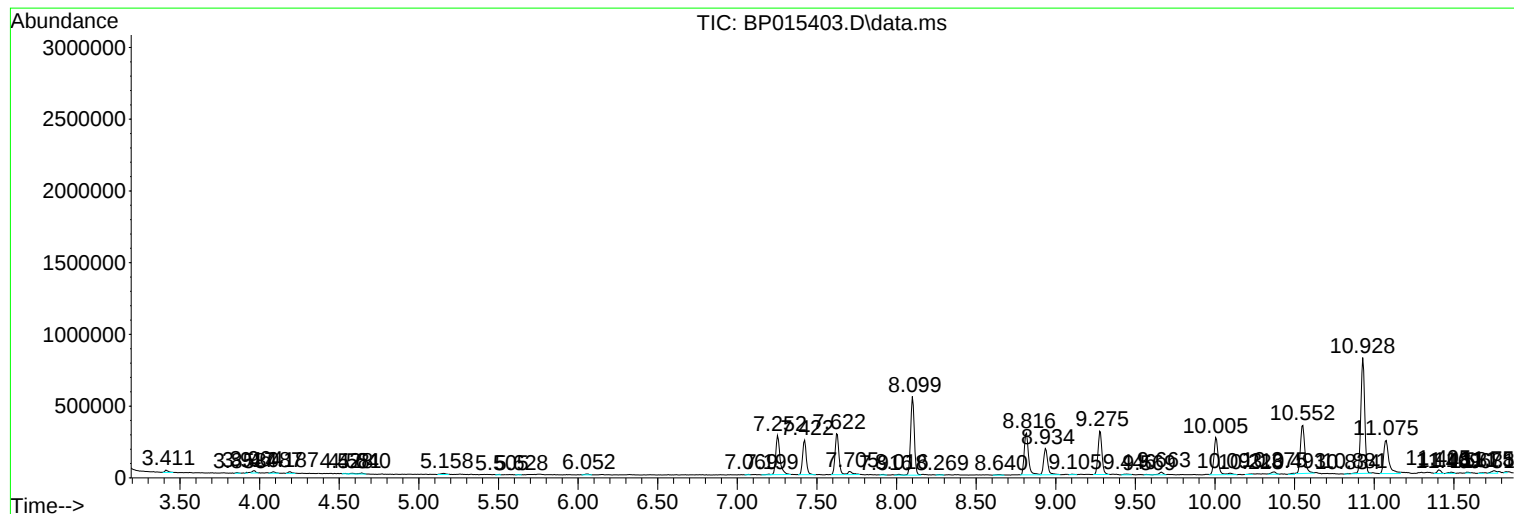
Sum of corrected areas: 38752377

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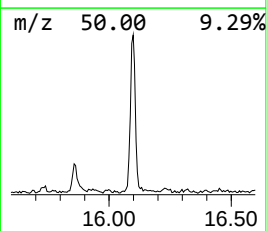
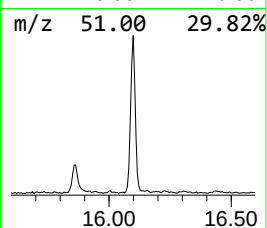
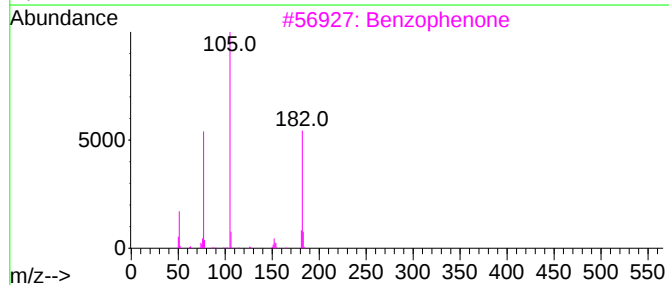
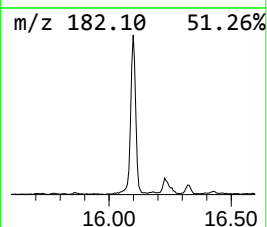
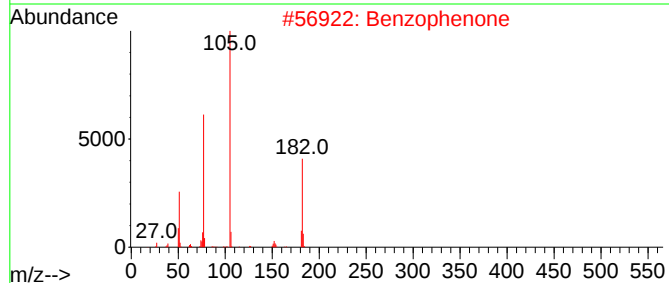
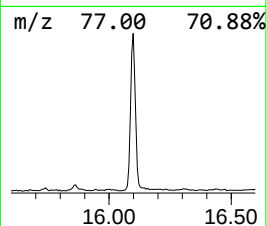
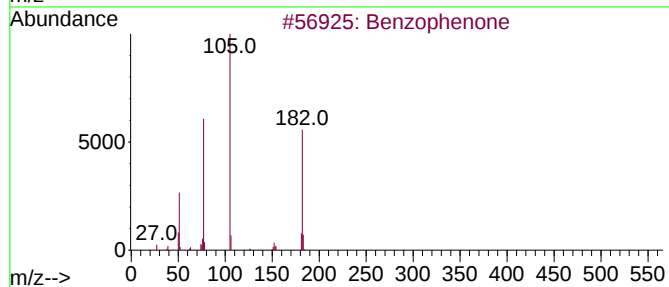
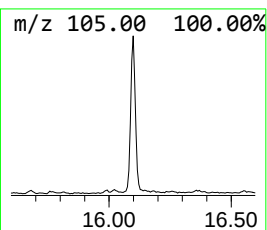
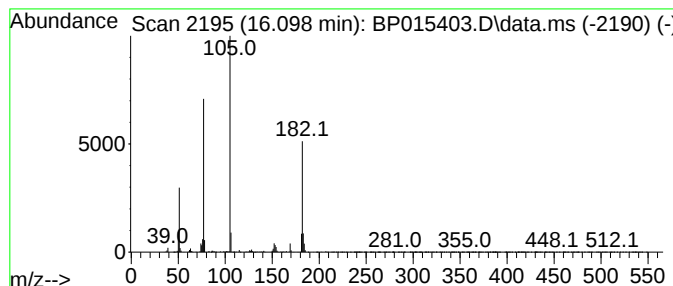
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.098	6.40 ng/ul	630603	Acenaphthene-d10	14.746

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	95
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	90
5			Benzophenone	182	C13H10O	000119-61-9	90



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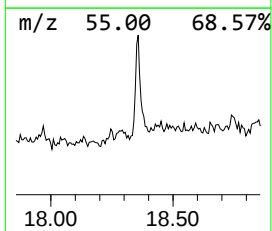
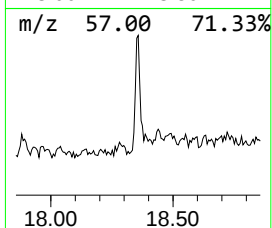
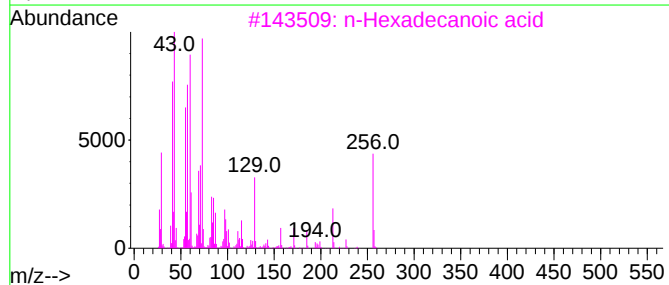
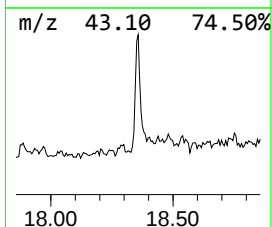
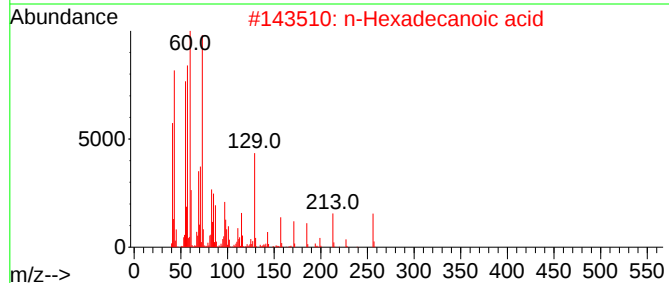
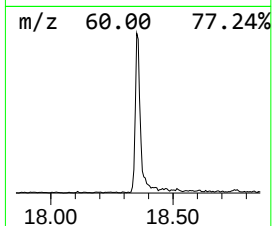
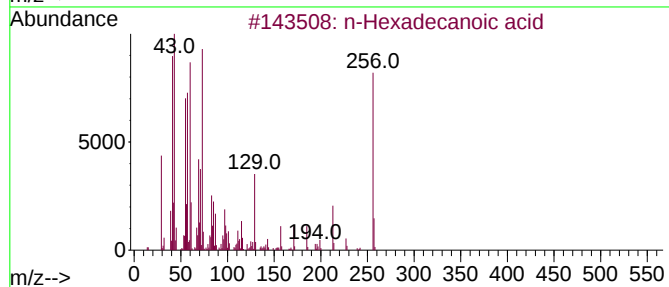
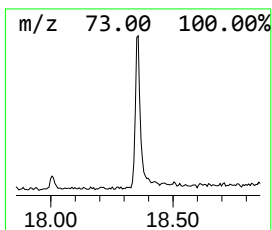
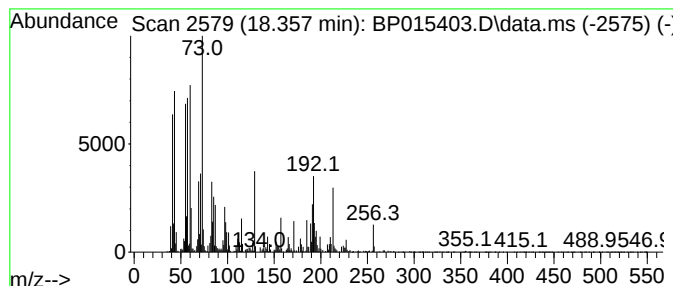
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.357	5.59 ng/ul	707645	Phenanthrene-d10	17.498

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	86
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	86



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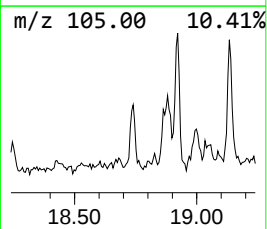
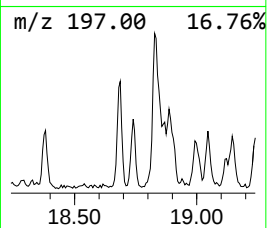
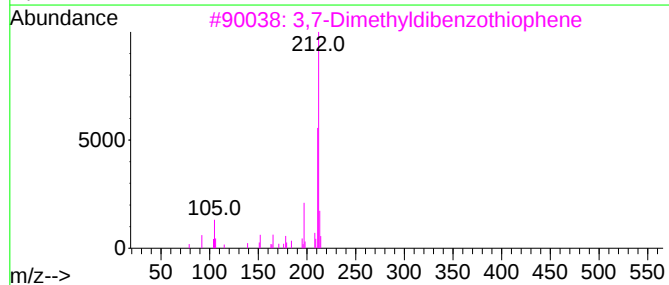
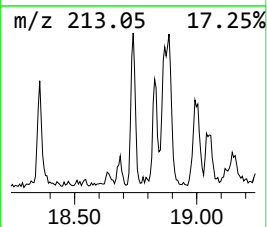
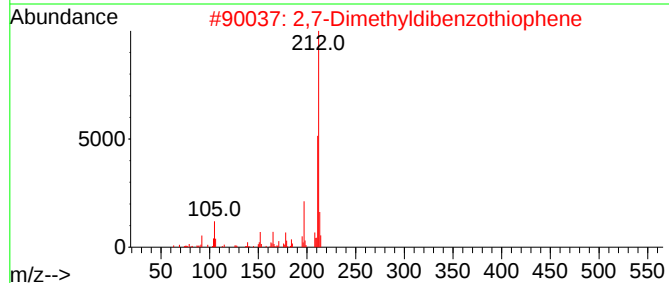
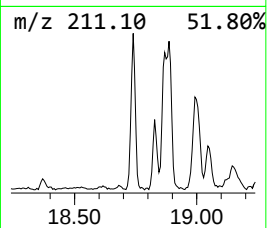
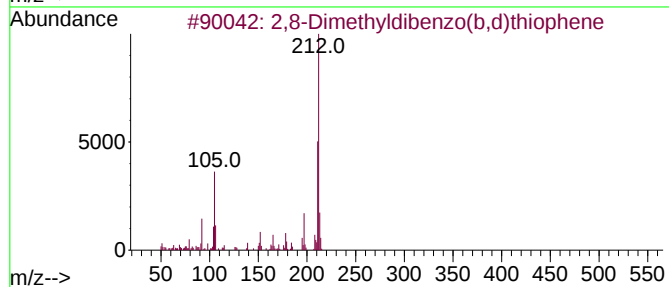
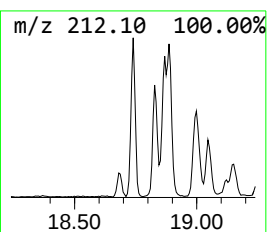
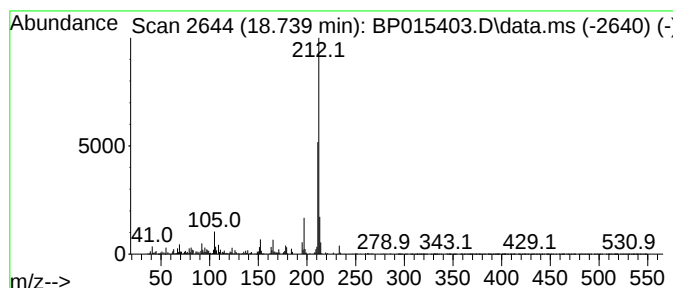
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2,8-Dimethyldibenzo(b,d)thi... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.739	2.45 ng/ul	310456	Phenanthrene-d10	17.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2,8-Dimethyldibenzo(b,d)thiophene	212	C14H12S	001207-15-4	94
2			2,7-Dimethyldibenzothiophene	212	C14H12S	031317-19-8	94
3			3,7-Dimethyldibenzothiophene	212	C14H12S	001136-85-2	93
4			2,6-Dimethyldibenzothiophene	212	C14H12S	089816-75-1	93
5			Dibenzothiophene, 4,6-dimethyl-	212	C14H12S	001207-12-1	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
Data File : BP015403.D
Acq On : 07 Jun 2023 17:34
Operator : MA/JU
Sample : 02950-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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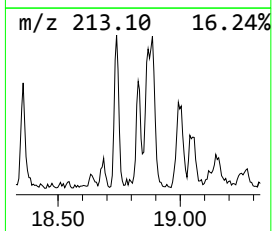
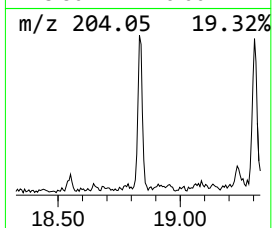
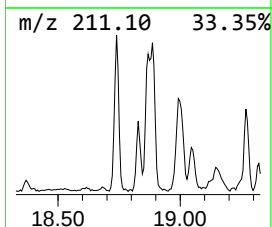
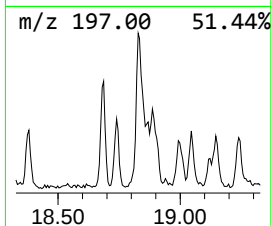
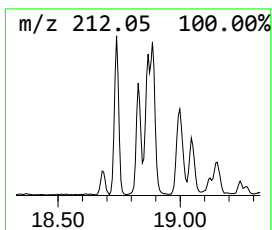
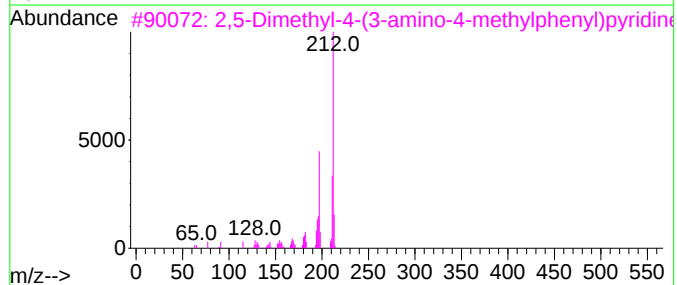
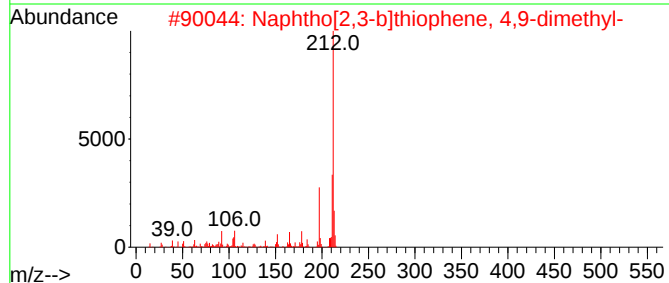
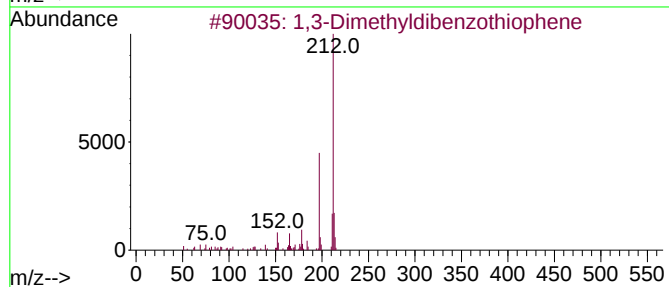
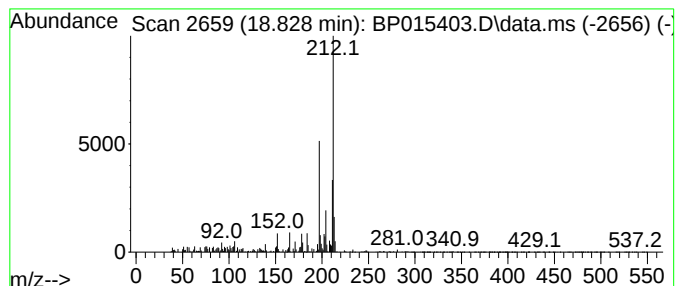
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1,3-Dimethyldibenzothiophene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.828	2.44 ng/ul	308945	Phenanthrene-d10	17.498

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1,3-Dimethyldibenzothiophene	212	C14H12S	1010383-55-0	90
2			Naphtho[2,3-b]thiophene, 4,9-dim...	212	C14H12S	016587-34-1	64
3			2,5-Dimethyl-4-(3-amino-4-methyl...	212	C14H16N2	071153-33-8	59
4			1,2-5,6-Dipyrzinecyclooctane	212	C12H12N4	099319-21-8	59
5			.beta.-Carboline, 8-methoxy-1-me...	212	C13H12N2O	212332-24-6	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
Data File : BP015403.D
Acq On : 07 Jun 2023 17:34
Operator : MA/JU
Sample : 02950-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW983

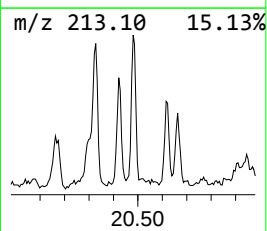
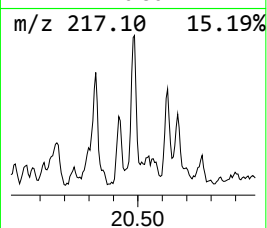
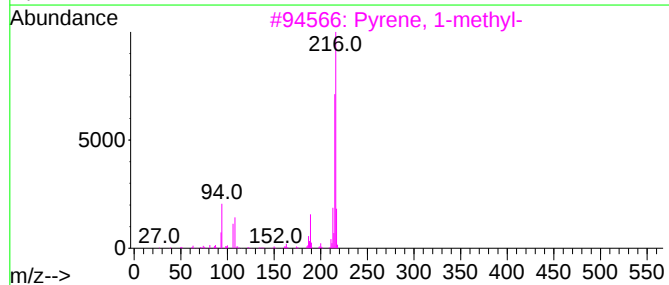
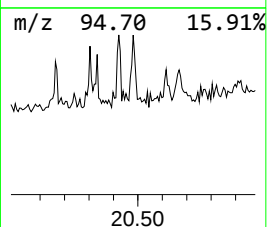
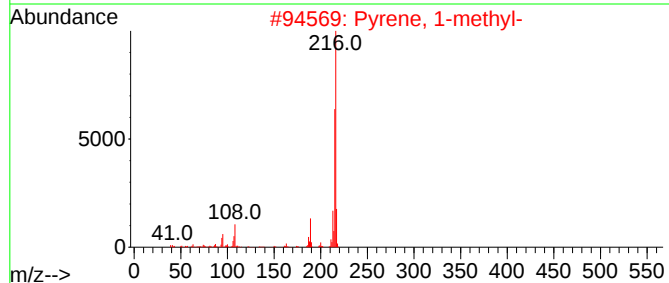
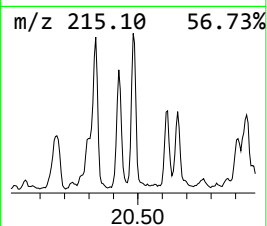
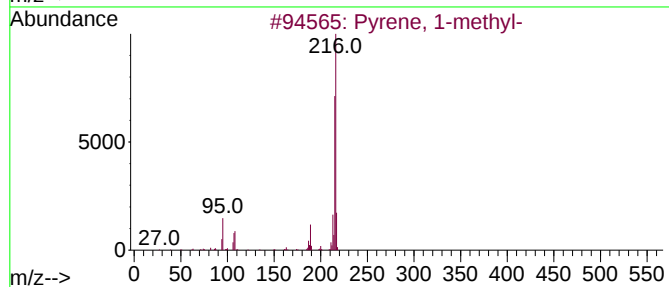
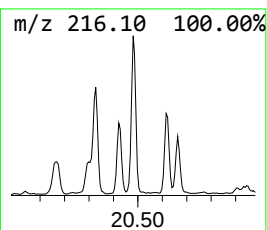
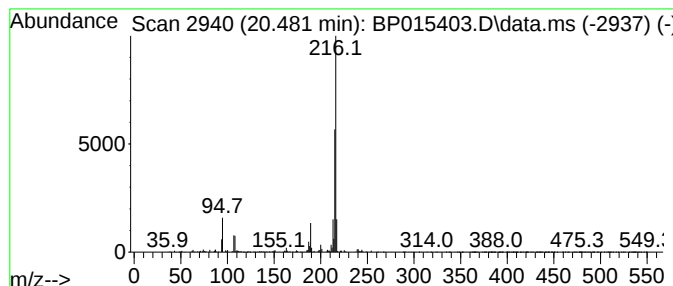
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Pyrene, 1-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
20.480	2.52 ng/ul	428233	Chrysene-d12	21.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pyrene, 1-methyl-	216	C17H12	002381-21-7	98
2			Pyrene, 1-methyl-	216	C17H12	002381-21-7	96
3			Pyrene, 1-methyl-	216	C17H12	002381-21-7	94
4			Pyrene, 2-methyl-	216	C17H12	003442-78-2	94
5			Pyrene, 4-methyl-	216	C17H12	003353-12-6	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
Data File : BP015403.D
Acq On : 07 Jun 2023 17:34
Operator : MA/JU
Sample : 02950-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EW983

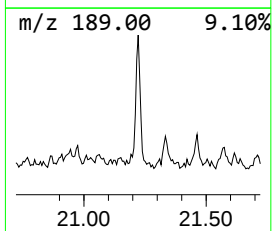
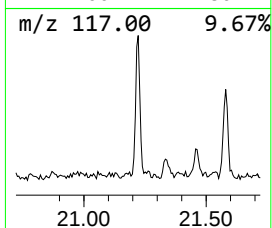
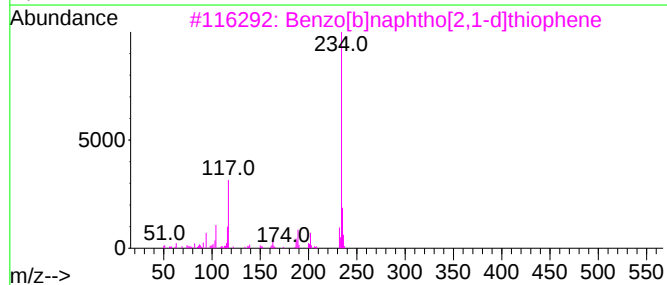
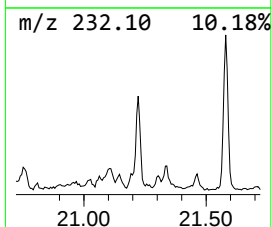
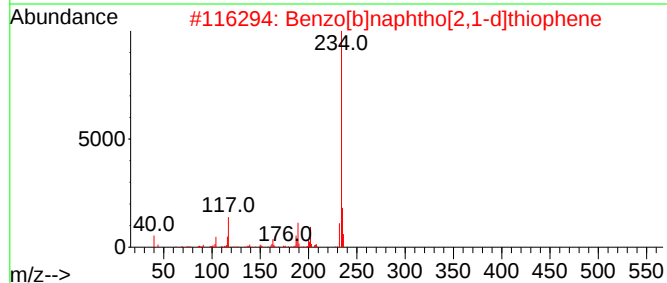
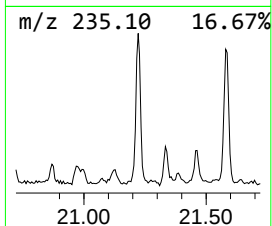
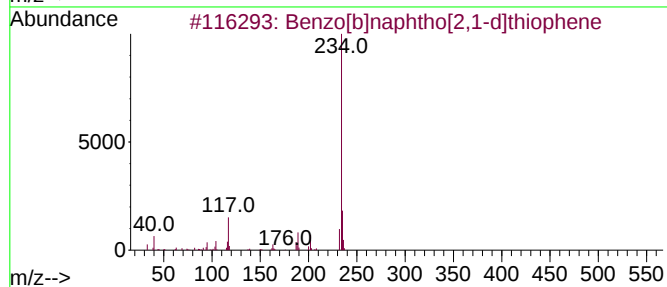
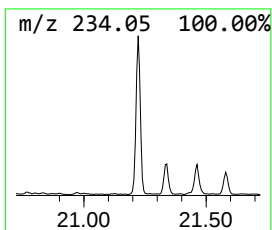
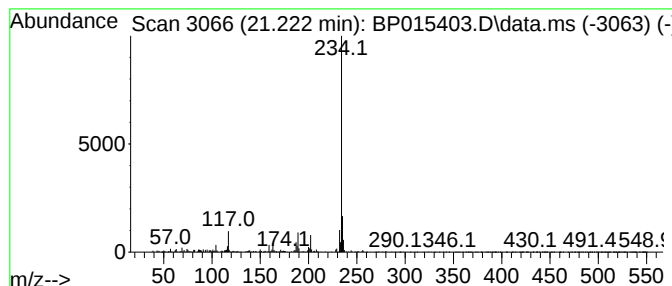
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzo[b]naphtho[2,1-d]thiop... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.222	3.01 ng/ul	511871	Chrysene-d12	21.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	97
2			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	96
3			Benzo[b]naphtho[2,1-d]thiophene	234	C16H10S	000239-35-0	94
4			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	94
5			Benzo[b]naphtho[2,3-d]thiophene	234	C16H10S	000243-46-9	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
Data File : BP015403.D
Acq On : 07 Jun 2023 17:34
Operator : MA/JU
Sample : 02950-06
Misc :
ALS Vial : 9 Sample Multiplier: 1

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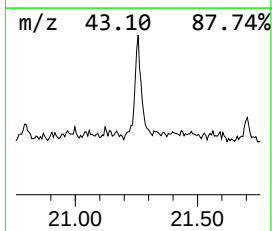
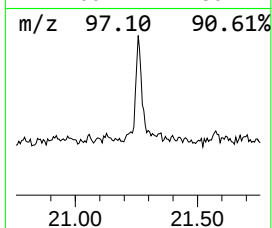
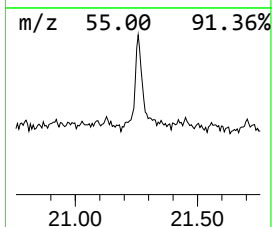
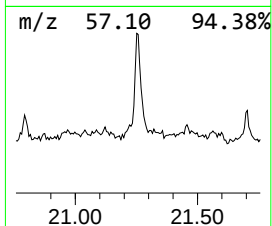
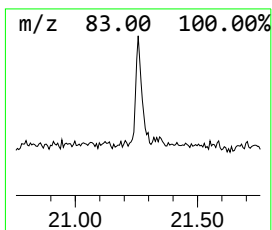
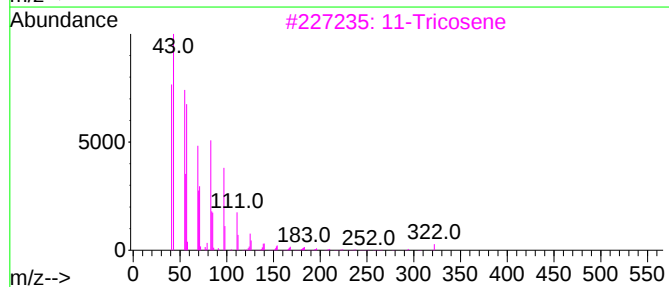
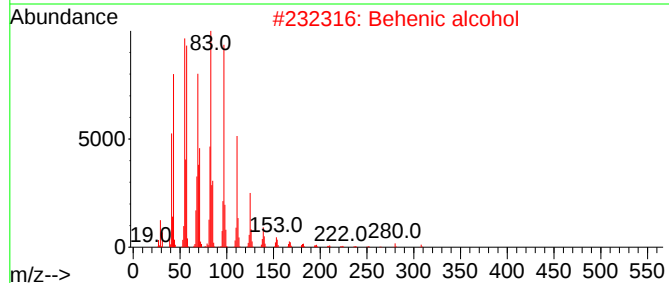
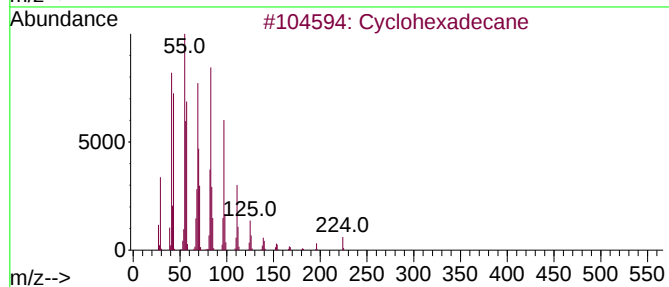
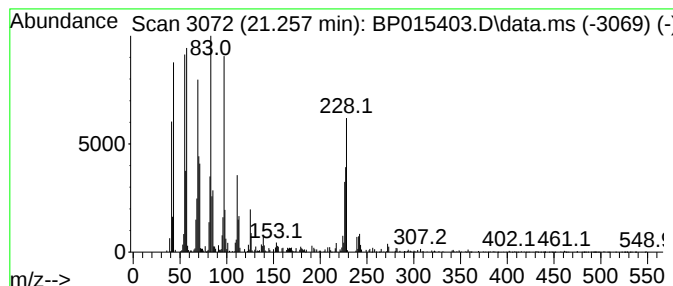
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP060723.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Cyclic21.257 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.257	3.34 ng/ul	567858	Chrysene-d12	21.581

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	91
2			Behenic alcohol	326	C22H46O	000661-19-8	89
3			11-Tricosene	322	C23H46	052078-56-5	76
4			1-Heneicosanol	312	C21H44O	015594-90-8	70
5			1-Nonadecene	266	C19H38	018435-45-5	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP060723\
Data File : BP015403.D
Acq On : 07 Jun 2023 17:34
Operator : MA/JU
Sample : 02950-06
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\NIST0.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	16.098	6.4	ng/ul	630603	3	14.746	1970550	20.0
n-Hexadecanoic ...	18.357	5.6	ng/ul	707645	4	17.498	2530160	20.0
2,8-Dimethyldib...	18.739	2.5	ng/ul	310456	4	17.498	2530160	20.0
1,3-Dimethyldib...	18.828	2.4	ng/ul	308945	4	17.498	2530160	20.0
Pyrene, 1-methyl-	20.480	2.5	ng/ul	428233	5	21.581	3404630	20.0
Benzo[b]naphtho...	21.222	3.0	ng/ul	511871	5	21.581	3404630	20.0
(DEL) Alkane: C...	21.257	3.3	ng/ul	567858	5	21.581	3404630	20.0