

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013257.D
 Acq On : 22 Dec 2022 22:26
 Operator : CG/JU
 Sample : N6015-18
 Misc :
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXY7

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-BP120722.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013257.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.334	21	25	31	rVB	215442	261392	1.76%	0.121%
2	3.758	94	97	124	rVB	2637468	3898724	26.25%	1.807%
3	4.087	150	153	159	rVB	55599	76716	0.52%	0.036%
4	5.040	307	315	321	rBV2	36498	68834	0.46%	0.032%
5	6.934	630	637	641	rBV4	27726	47868	0.32%	0.022%
6	7.111	659	667	678	rBV	3466375	5635088	37.95%	2.611%
7	7.275	688	695	707	rVB	2796411	4290575	28.89%	1.988%
8	7.475	722	729	740	rBV	3600294	5599405	37.71%	2.595%
9	7.946	800	809	816	rBV	2283126	3550696	23.91%	1.645%
10	8.657	923	930	947	rBV	4233677	6779372	45.65%	3.141%
11	9.116	1000	1008	1015	rBV	3368932	5402222	36.38%	2.503%
12	9.516	1071	1076	1081	rBV2	55397	88702	0.60%	0.041%
13	9.840	1122	1131	1141	rBV	2832536	4707163	31.70%	2.181%
14	10.381	1214	1223	1234	rBV	4273241	7157273	48.20%	3.316%
15	10.757	1276	1287	1293	rBV	2968588	5009030	33.73%	2.321%
16	10.810	1293	1296	1303	rVB	211573	317379	2.14%	0.147%
17	10.899	1303	1311	1330	rBV	3571928	6026560	40.58%	2.793%
18	11.769	1454	1459	1467	rVB4	26396	52724	0.36%	0.024%
19	12.010	1496	1500	1506	rVB2	16480	27031	0.18%	0.013%
20	12.075	1506	1511	1516	rBV	23164	40742	0.27%	0.019%
21	12.169	1521	1527	1533	rVB2	44960	73331	0.49%	0.034%
22	12.346	1552	1557	1563	rVV2	69483	113778	0.77%	0.053%
23	12.416	1563	1569	1578	rVB	75829	126639	0.85%	0.059%
24	12.546	1583	1591	1599	rBV	10412	28976	0.20%	0.013%
25	12.634	1599	1606	1611	rVB	39954	62679	0.42%	0.029%
26	12.851	1637	1643	1649	rBV	15138	36148	0.24%	0.017%
27	13.110	1681	1687	1694	rBV4	32651	56473	0.38%	0.026%
28	13.404	1728	1737	1744	rBV2	92130	182890	1.23%	0.085%
29	13.575	1758	1766	1770	rVB4	32927	70762	0.48%	0.033%
30	13.763	1789	1798	1805	rBV7	21545	65893	0.44%	0.031%
31	13.904	1817	1822	1828	rBV2	38745	72638	0.49%	0.034%
32	13.993	1829	1837	1844	rVV	6206744	8993147	60.56%	4.167%
33	14.057	1846	1848	1852	rVV2	41742	46004	0.31%	0.021%
34	14.110	1852	1857	1861	rVV6	33611	51195	0.34%	0.024%
35	14.175	1865	1868	1872	rVB	49976	59587	0.40%	0.028%
36	14.228	1872	1877	1880	rBV5	19795	38280	0.26%	0.018%

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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-BP120722.M
 Title : SVOA CALIBRATION

37	14.281	1880	1886	1897	rVV	7743992	11500146	77.44%	5.329%
38	14.475	1915	1919	1923	rVB	68897	91925	0.62%	0.043%
39	14.587	1928	1938	1945	rVV	4125363	5950660	40.07%	2.757%
40	14.651	1945	1949	1957	rVB2	93690	164853	1.11%	0.076%
41	14.787	1966	1972	1987	rBV	2221326	3592285	24.19%	1.665%
42	14.945	1995	1999	2001	rBV4	23136	30349	0.20%	0.014%
43	14.987	2002	2006	2013	rVB	119787	184945	1.25%	0.086%
44	15.204	2038	2043	2047	rBV6	20785	44310	0.30%	0.021%
45	15.375	2067	2072	2075	rBV6	27677	42842	0.29%	0.020%
46	15.422	2077	2080	2084	rVV	38467	50196	0.34%	0.023%
47	15.581	2099	2107	2114	rBV	9043177	12913236	86.96%	5.984%
48	15.634	2114	2116	2121	rVV	78293	97505	0.66%	0.045%
49	15.704	2122	2128	2136	rVV	2304702	3247700	21.87%	1.505%
50	15.822	2144	2148	2156	rVB5	48654	98626	0.66%	0.046%
51	15.945	2163	2169	2182	rVB	3806443	5152845	34.70%	2.388%
52	16.081	2188	2192	2198	rVB2	36314	71157	0.48%	0.033%
53	16.157	2198	2205	2206	rBV7	16295	31706	0.21%	0.015%
54	16.292	2224	2228	2230	rBV2	74058	112591	0.76%	0.052%
55	16.516	2261	2266	2273	rVB	156464	206592	1.39%	0.096%
56	16.728	2299	2302	2308	rBV5	40967	64547	0.43%	0.030%
57	16.875	2324	2327	2331	rBV3	23936	27543	0.19%	0.013%
58	16.916	2331	2334	2338	rVB5	24345	30213	0.20%	0.014%
59	16.998	2343	2348	2355	rBV2	84721	148811	1.00%	0.069%
60	17.092	2358	2364	2368	rBV3	66424	113965	0.77%	0.053%
61	17.163	2373	2376	2380	rVB	55681	75234	0.51%	0.035%
62	17.228	2383	2387	2391	rBV	80395	92604	0.62%	0.043%
63	17.292	2395	2398	2401	rBV4	32665	41725	0.28%	0.019%
64	17.345	2401	2407	2411	rBV	4104694	5788236	38.98%	2.682%
65	17.386	2411	2414	2418	rVV	1172943	1575760	10.61%	0.730%
66	17.445	2418	2424	2436	rVB	8525780	12286866	82.74%	5.693%
67	17.698	2464	2467	2469	rBV2	55165	62318	0.42%	0.029%
68	17.745	2473	2475	2481	rVB2	75551	106617	0.72%	0.049%
69	17.863	2492	2495	2500	rVB4	58535	73647	0.50%	0.034%
70	18.004	2515	2519	2521	rBV5	43854	56414	0.38%	0.026%
71	18.104	2531	2536	2539	rBV5	63299	99308	0.67%	0.046%
72	18.157	2542	2545	2549	rBV3	81659	102116	0.69%	0.047%
73	18.216	2550	2555	2559	rBV2	1483201	1976796	13.31%	0.916%
74	18.257	2559	2562	2569	rVB	256007	386041	2.60%	0.179%
75	18.398	2581	2586	2590	rBV2	215947	384122	2.59%	0.178%
76	18.498	2600	2603	2608	rBV3	84347	132779	0.89%	0.062%

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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

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77	18.692	2632	2636	2639	rVB	138282	169603	1.14%	0.079%
78	18.733	2639	2643	2648	rVB2	95955	127913	0.86%	0.059%
79	19.098	2700	2705	2710	rBV2	144217	307881	2.07%	0.143%
80	19.351	2738	2748	2754	rBV	3060184	4324118	29.12%	2.004%
81	19.445	2761	2764	2769	rBV2	282313	335339	2.26%	0.155%
82	19.498	2770	2773	2777	rVB	149691	186670	1.26%	0.086%
83	19.586	2786	2788	2796	rVB3	103519	174663	1.18%	0.081%
84	19.675	2798	2803	2807	rVB	10004882	14850135	100.00%	6.881%
85	19.875	2835	2837	2842	rVB	115606	131297	0.88%	0.061%
86	20.016	2857	2861	2868	rBV2	156071	378959	2.55%	0.176%
87	20.186	2887	2890	2896	rVB	451692	597338	4.02%	0.277%
88	20.280	2903	2906	2910	rBV2	196878	257546	1.73%	0.119%
89	20.480	2936	2940	2942	rBV	186527	264678	1.78%	0.123%
90	20.916	3011	3014	3020	rBV	270834	381607	2.57%	0.177%
91	21.127	3045	3050	3054	rBV2	1862922	2797263	18.84%	1.296%
92	21.433	3095	3102	3106	rBV2	5392233	9866611	66.44%	4.572%
93	21.469	3106	3108	3115	rVB	1578191	2116356	14.25%	0.981%
94	23.151	3386	3394	3400	rBV2	1674354	4988808	33.59%	2.312%
95	23.698	3482	3487	3490	rBV	868701	1607787	10.83%	0.745%
96	23.751	3490	3496	3502	rBV2	7472256	14654877	98.69%	6.791%
97	23.916	3517	3524	3530	rBV2	3229895	6753531	45.48%	3.129%
98	24.421	3604	3610	3621	rVB4	1546637	3448505	23.22%	1.598%
99	24.527	3624	3628	3637	rVB	1077294	2230082	15.02%	1.033%
100	29.003	4380	4389	4402	rVB4	2272837	8532490	57.46%	3.954%

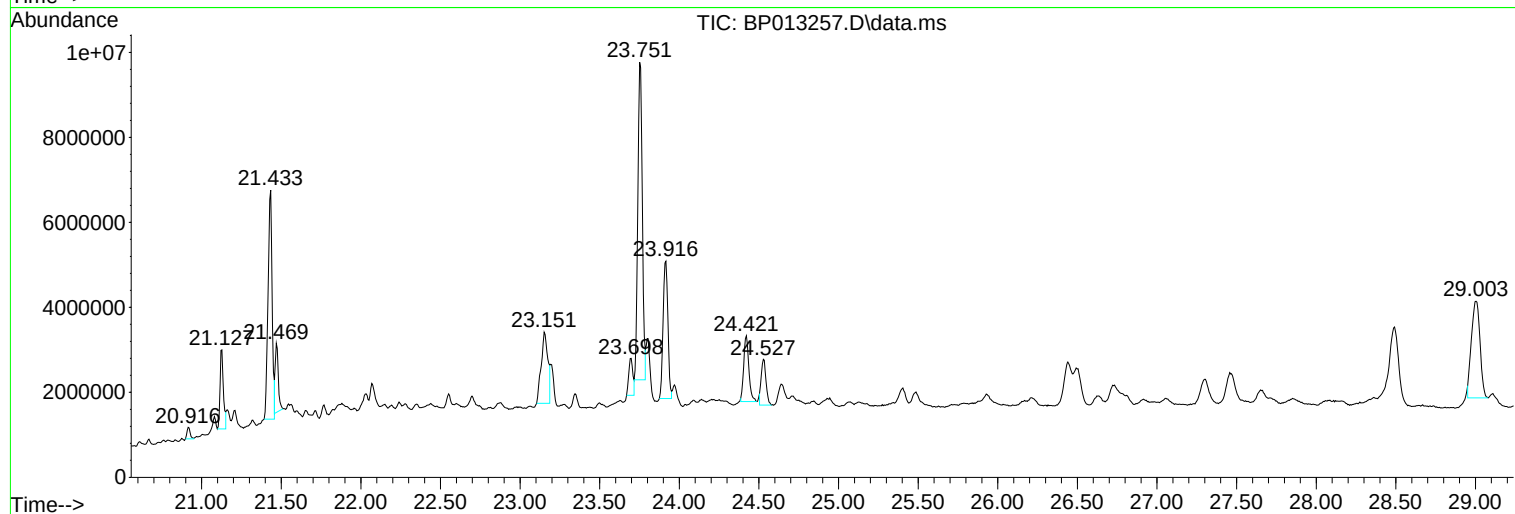
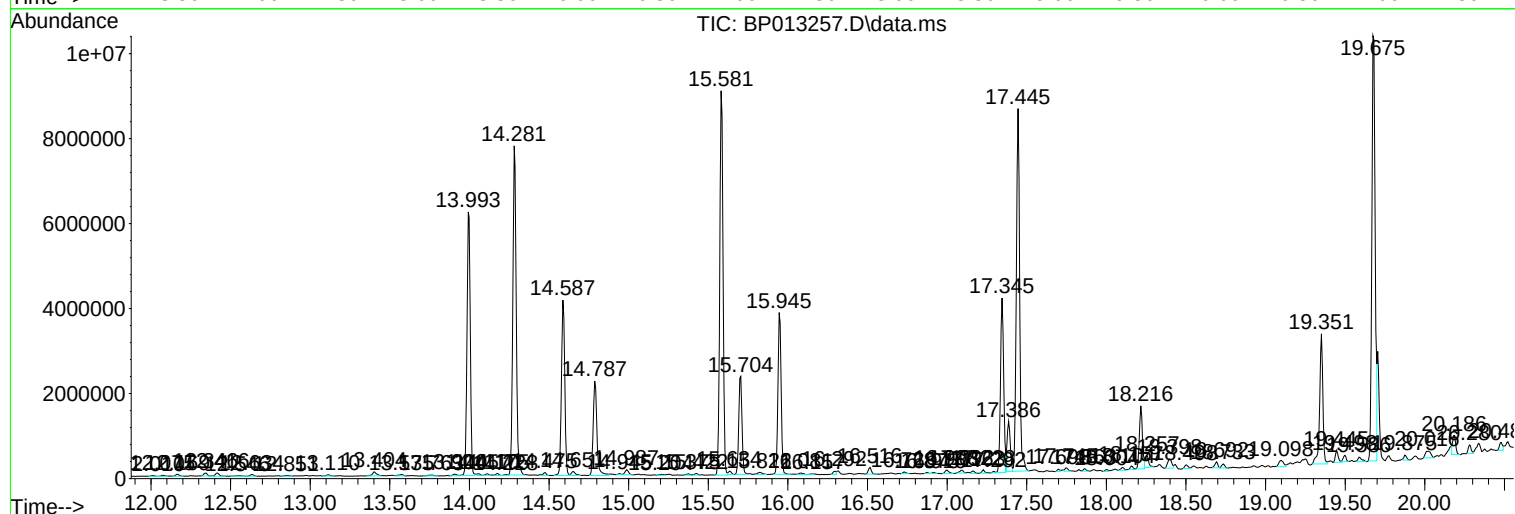
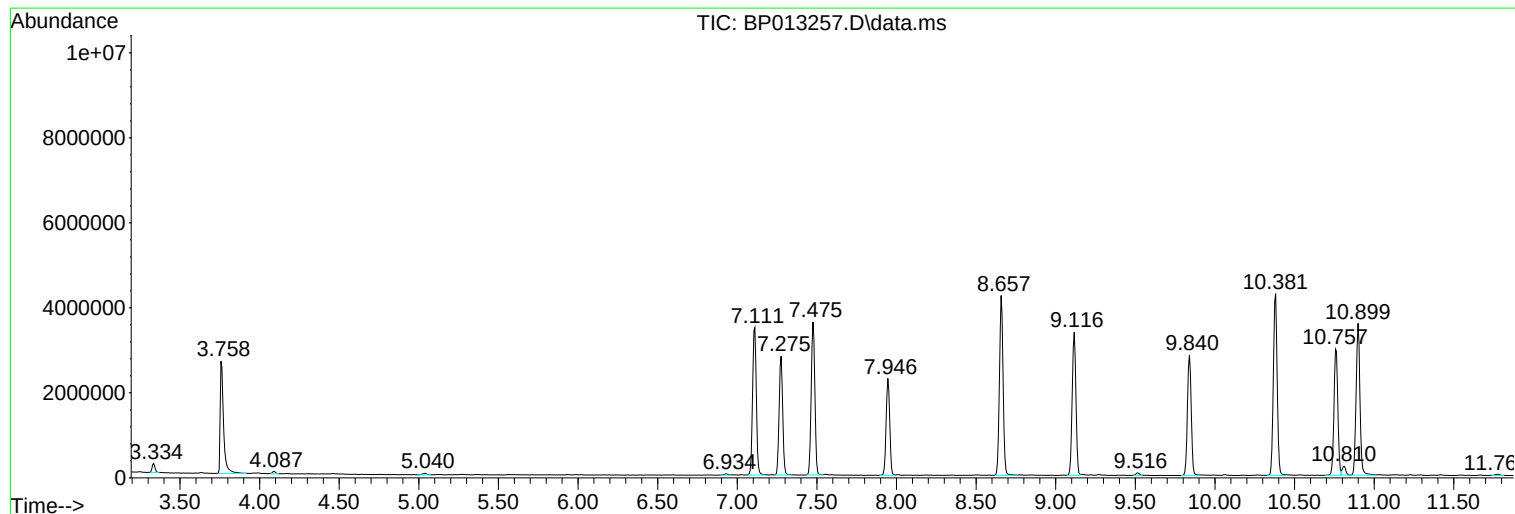
Sum of corrected areas: 215810104

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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



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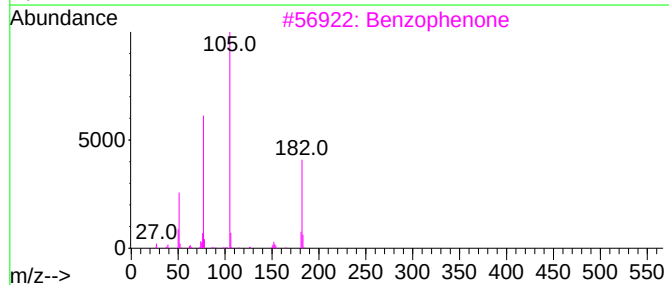
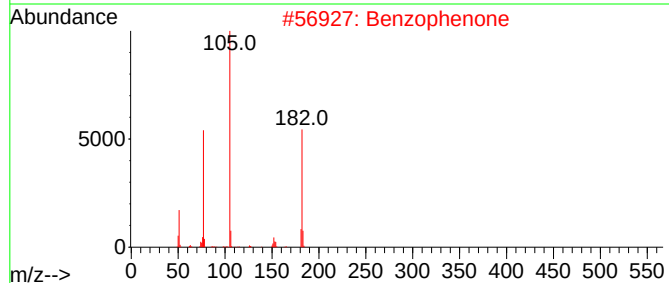
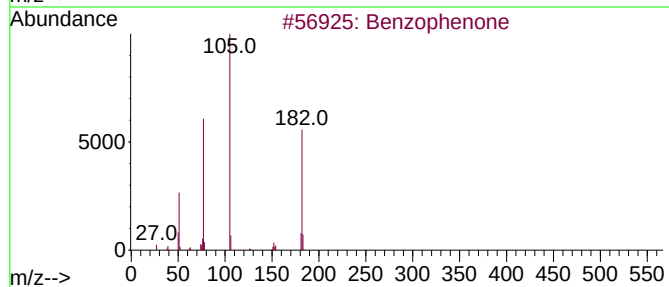
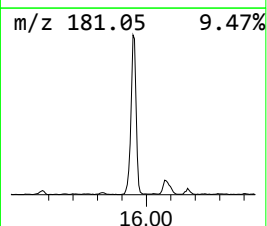
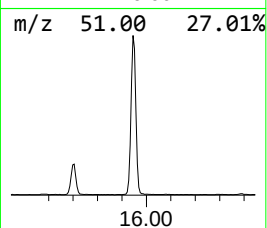
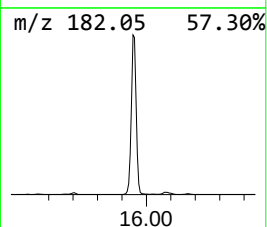
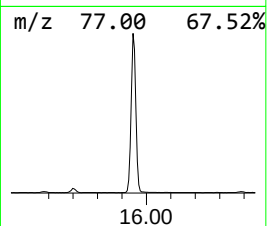
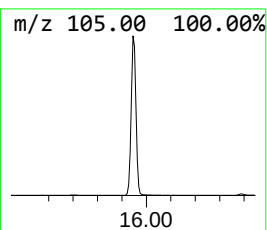
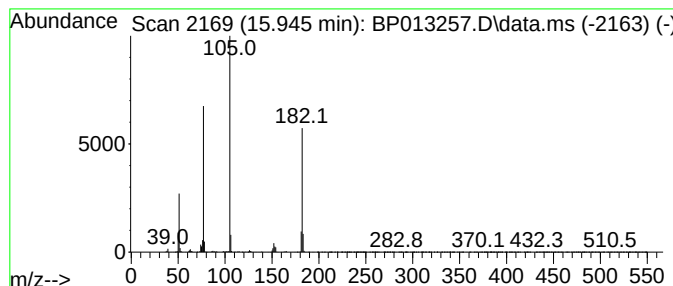
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	17.32 ng/ul	5152850	Acenaphthene-d10	14.587

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	96
2			Benzophenone	182	C13H10O	000119-61-9	95
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



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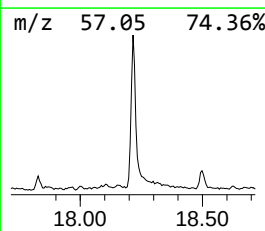
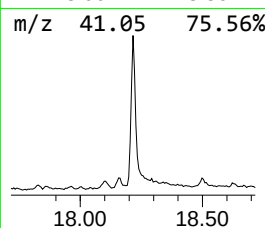
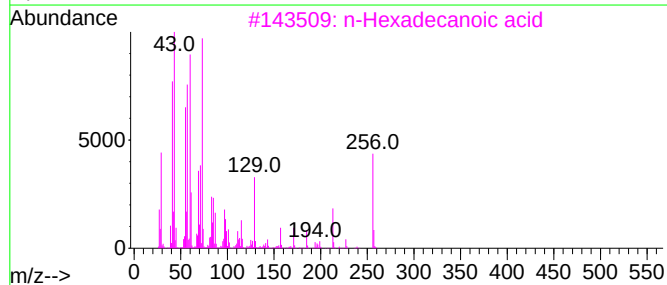
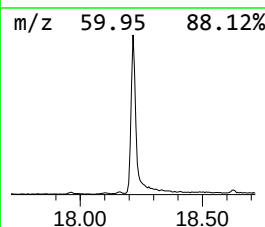
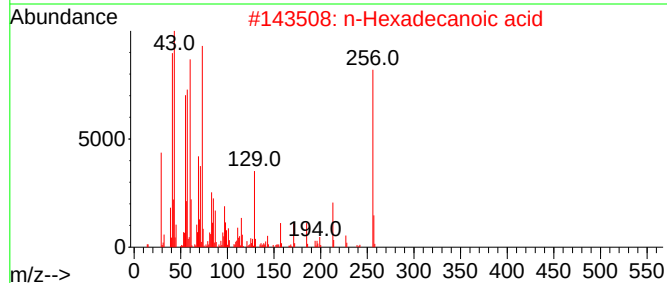
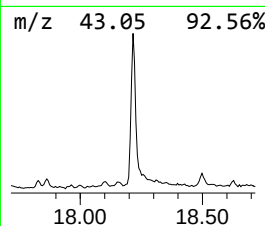
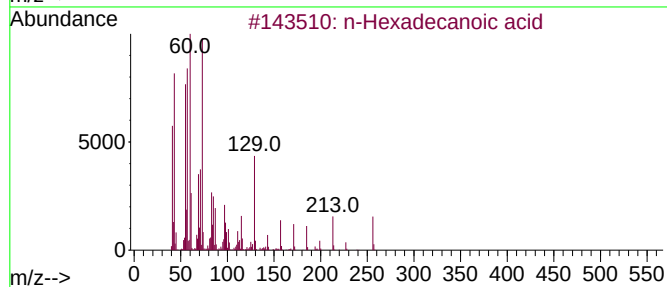
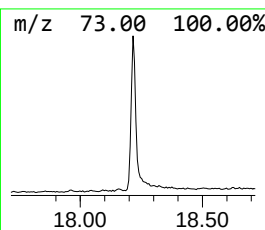
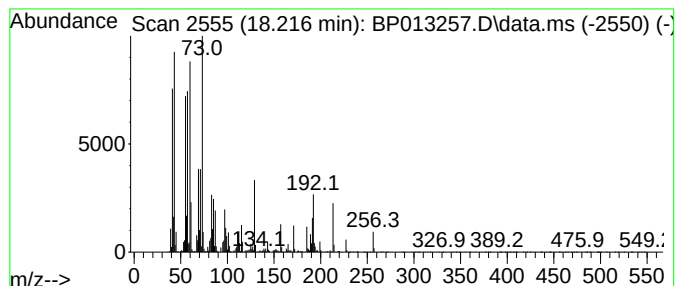
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	6.83 ng/ul	1976800	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
2			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	99
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	93
5			Tetradecanoic acid	228	C14H28O2	000544-63-8	89



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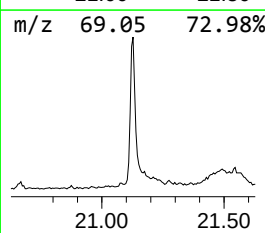
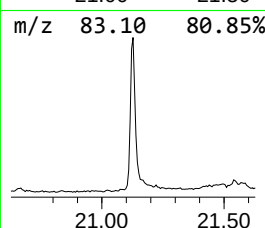
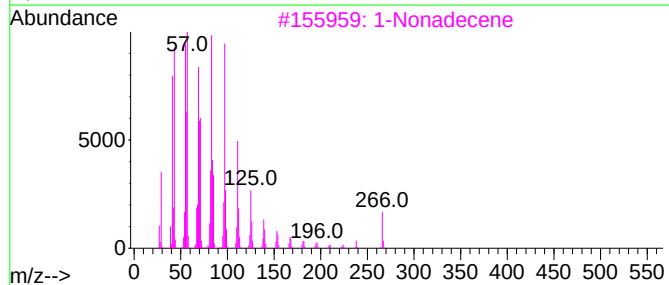
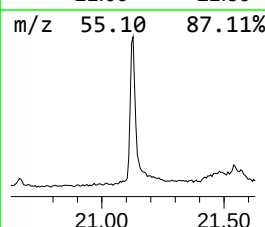
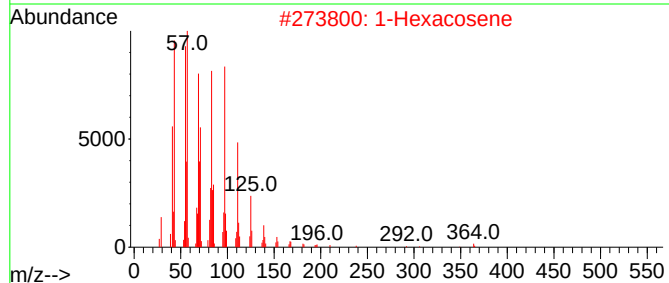
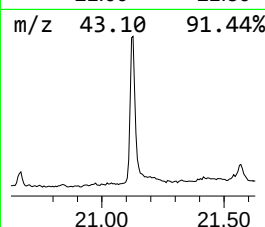
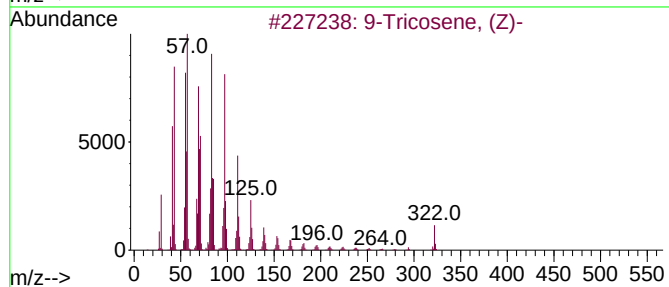
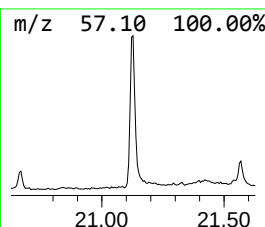
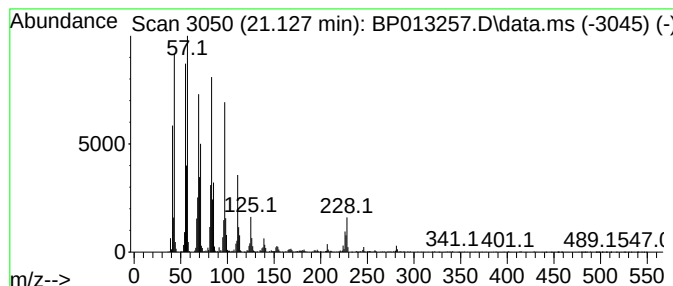
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	5.67 ng/ul	2797260	Chrysene-d12	21.433

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Tricosene, (Z)-		322	C23H46	027519-02-4	99
2	1-Hexacosene		364	C26H52	018835-33-1	99
3	1-Nonadecene		266	C19H38	018435-45-5	98
4	5-Eicosene, (E)-		280	C20H40	074685-30-6	97
5	1-Heneicosyl formate		340	C22H44O2	077899-03-7	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013257.D
Acq On : 22 Dec 2022 22:26
Operator : CG/JU
Sample : N6015-18
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXY7

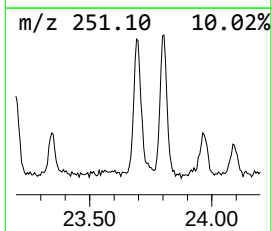
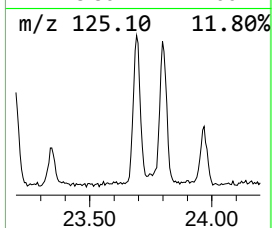
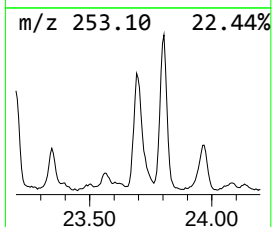
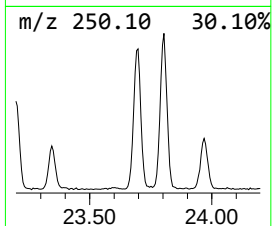
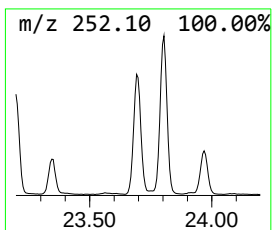
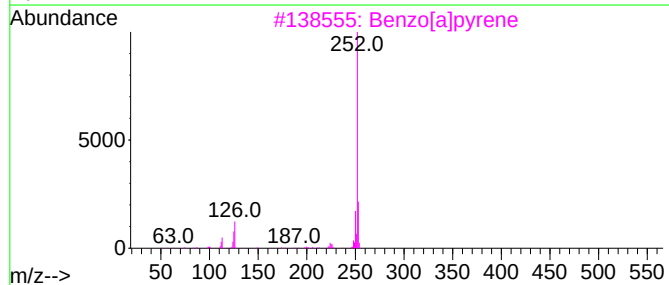
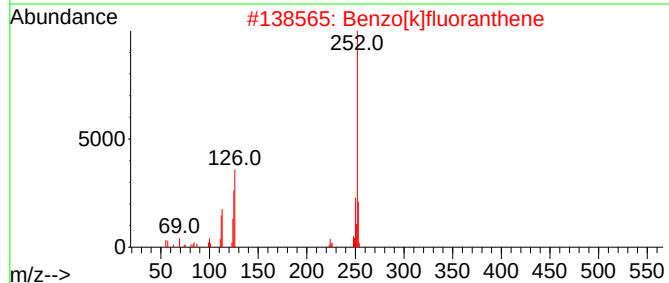
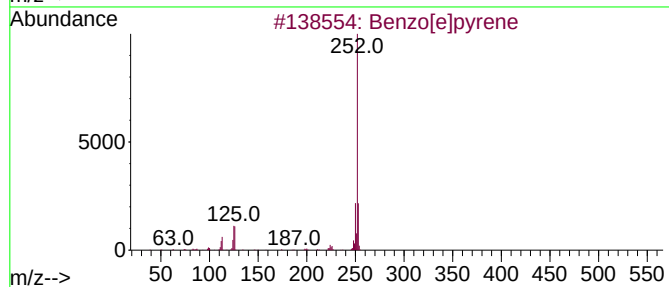
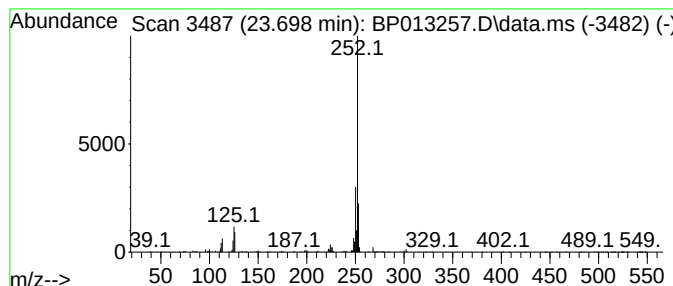
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.698	4.76 ng/ul	1607790	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	98
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	94
3			Benzo[a]pyrene	252	C20H12	000050-32-8	93
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93
5			Benzo[b]fluoranthene	252	C20H12	000205-99-2	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013257.D
Acq On : 22 Dec 2022 22:26
Operator : CG/JU
Sample : N6015-18
Misc :
ALS Vial : 24 Sample Multiplier: 1

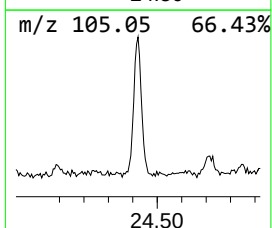
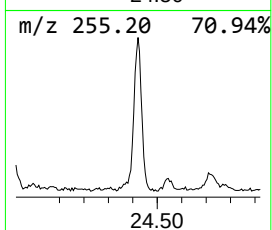
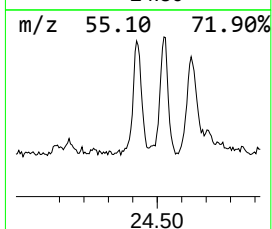
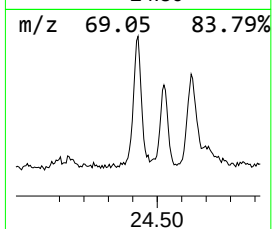
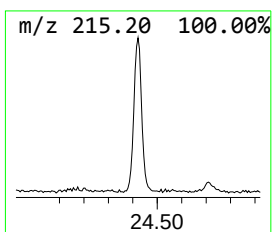
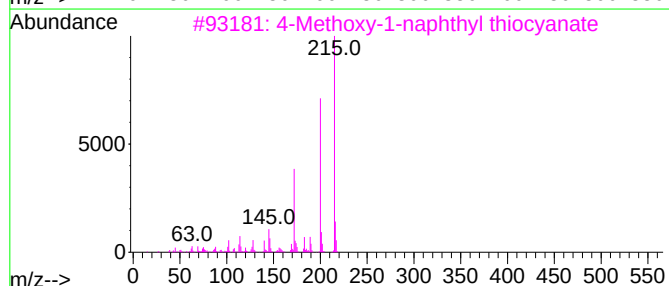
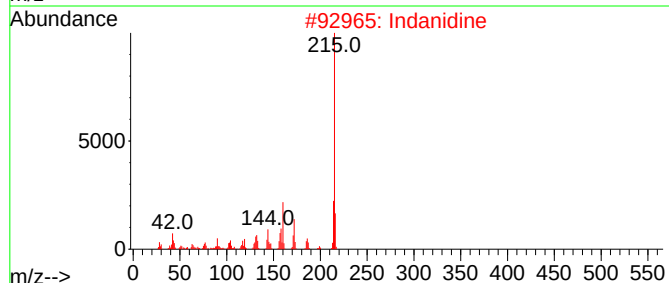
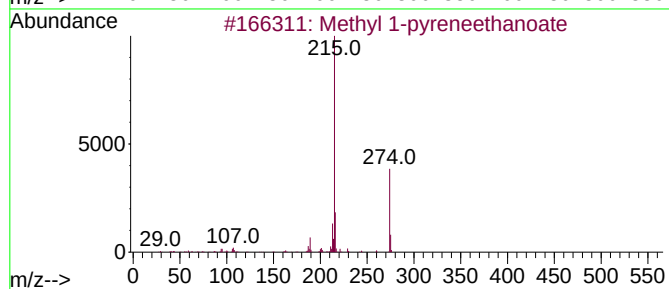
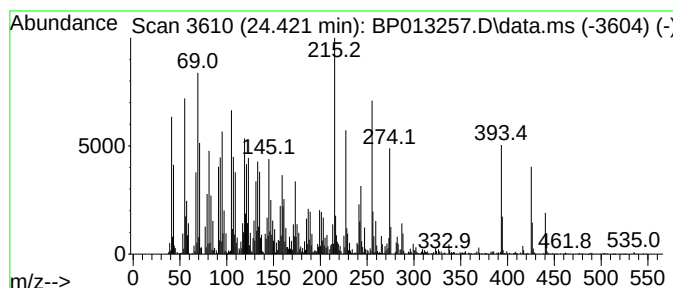
Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-01 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.421	10.21 ng/ul	3448510	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	43
2	Indanidine	215	C11H13N5	085392-79-6	15
3	4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	11
4	N-[2-(1H-Indol-3-yl)ethyl]propan...	288	C16H24N2OSi	1000493-19-0	11
5	Ethyl 4-[3-trifluoromethylphenyl]...	288	C13H11F3O4	1000211-86-0	11



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013257.D
Acq On : 22 Dec 2022 22:26
Operator : CG/JU
Sample : N6015-18
Misc :
ALS Vial : 24 Sample Multiplier: 1

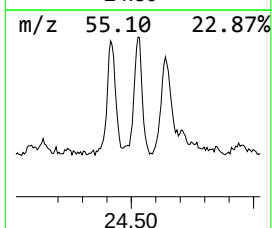
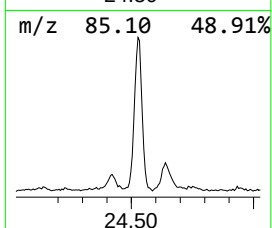
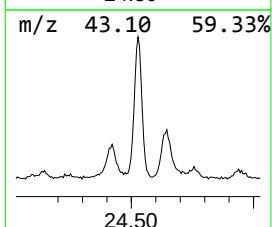
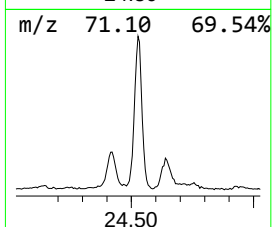
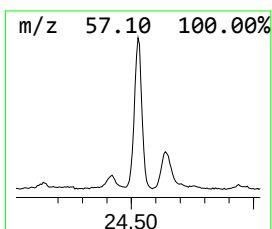
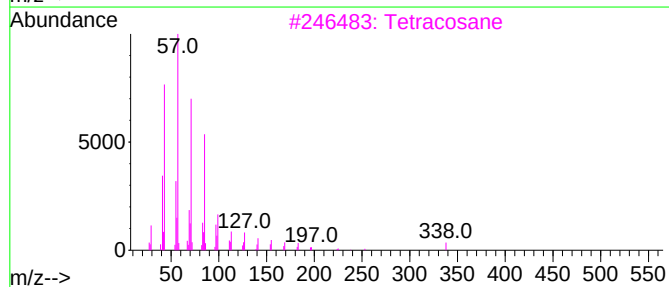
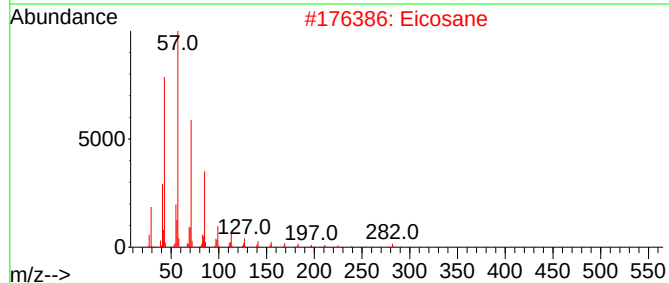
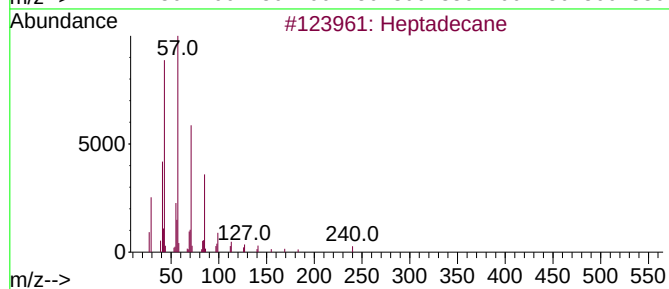
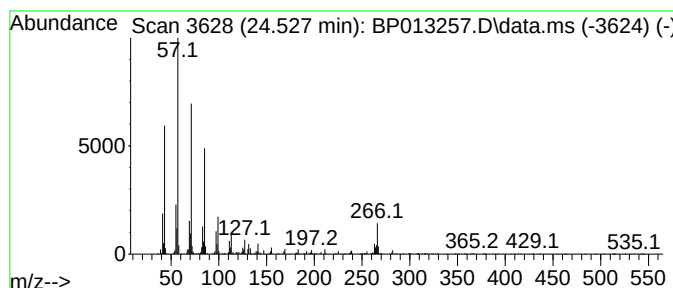
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
24.527	6.60 ng/ul	2230080	Perylene-d12	23.916	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Heptadecane	240	C17H36	000629-78-7	94
2	Eicosane	282	C20H42	000112-95-8	93
3	Tetracosane	338	C24H50	000646-31-1	93
4	Hexacosane	366	C26H54	000630-01-3	87
5	Nonacosane, 2-methyl-	422	C30H62	001560-75-4	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013257.D
Acq On : 22 Dec 2022 22:26
Operator : CG/JU
Sample : N6015-18
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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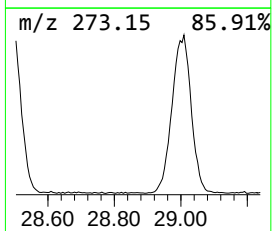
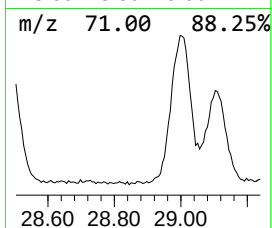
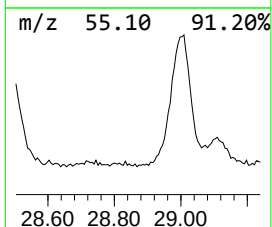
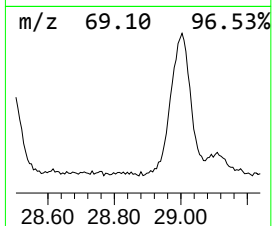
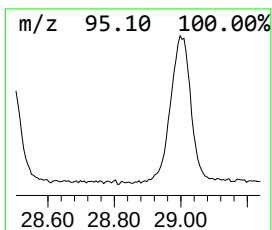
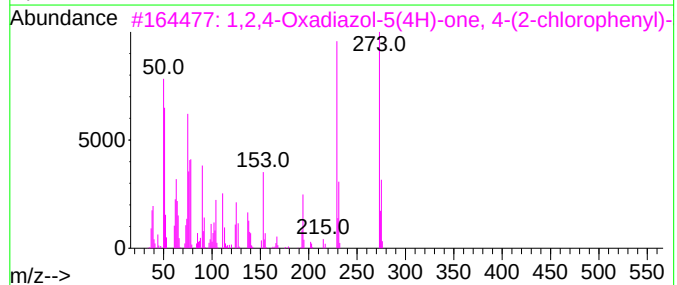
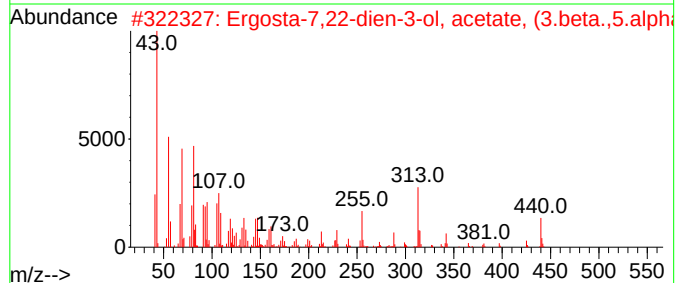
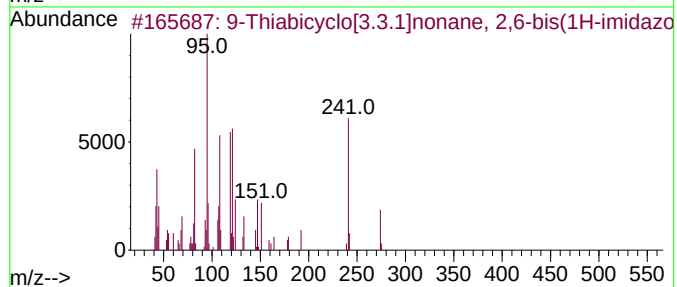
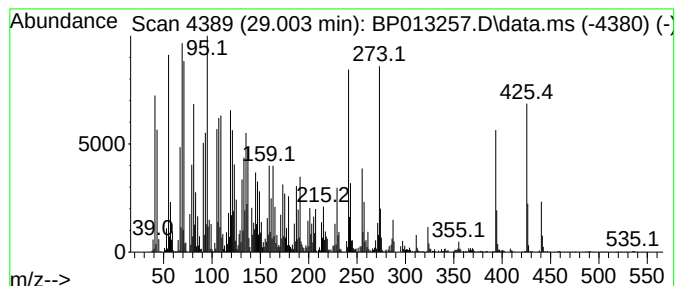
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.003	25.27 ng/ul	8532490	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Thiabicyclo[3.3.1]nonane, 2,6-...	274	C14H18N4S	1000141-52-1	30		
2	Ergosta-7,22-dien-3-ol, acetate,...	440	C30H48O2	001449-60-1	25		
3	1,2,4-Oxadiazol-5(4H)-one, 4-(2-...	273	C13H8ClN3O2	288246-55-9	25		
4	Silane, dimethyl(2-chlorophenoxy...	440	C25H45ClO2Si	1000347-08-2	25		
5	Malonodinitrile, 2-[3-dimethylam...	273	C18H15N3	1000264-78-5	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
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Instrument :
BNA_P
ClientSampleId :
DBXY7

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.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	15.945	17.3	ng/u1	5152850	3	14.587	5950660	20.0
n-Hexadecanoic ...	18.216	6.8	ng/u1	1976800	4	17.345	5788240	20.0
(DEL) Alkane: S...	21.127	5.7	ng/u1	2797260	5	21.433	9866610	20.0
Benzo[e]pyrene	23.698	4.8	ng/u1	1607790	6	23.916	6753530	20.0
unknown-01	24.421	10.2	ng/u1	3448510	6	23.916	6753530	20.0
(DEL) Alkane: S...	24.527	6.6	ng/u1	2230080	6	23.916	6753530	20.0
unknown-02	29.003	25.3	ng/u1	8532490	6	23.916	6753530	20.0