

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013256.D
 Acq On : 22 Dec 2022 21:51
 Operator : CG/JU
 Sample : N6015-17
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXY5

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: BP013256.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	32	rVB	187363	231141	1.33%	0.088%
2	3.758	94	97	113	rBV	2471975	3731491	21.49%	1.423%
3	4.087	149	153	158	rBV	55680	77658	0.45%	0.030%
4	5.040	308	315	321	rVB2	42042	74446	0.43%	0.028%
5	7.105	658	666	679	rBV	3314895	5332565	30.71%	2.034%
6	7.275	688	695	708	rVB	2651104	4117367	23.71%	1.571%
7	7.475	720	729	736	rBV	3441416	5326234	30.67%	2.032%
8	7.946	799	809	816	rBV	2429869	3741856	21.55%	1.427%
9	8.658	922	930	941	rBV	4060190	6421592	36.98%	2.449%
10	9.116	1000	1008	1015	rVV	3209654	5208644	30.00%	1.987%
11	9.275	1029	1035	1040	rVB4	30712	51848	0.30%	0.020%
12	9.516	1069	1076	1084	rVV	74439	129080	0.74%	0.049%
13	9.840	1123	1131	1143	rBV	2840800	4593219	26.45%	1.752%
14	10.381	1215	1223	1234	rBV	4270647	7146263	41.15%	2.726%
15	10.757	1276	1287	1293	rBV	3280583	5539833	31.90%	2.113%
16	10.810	1293	1296	1304	rVV	298640	462102	2.66%	0.176%
17	10.899	1304	1311	1330	rVV	3348056	5881679	33.87%	2.244%
18	12.163	1521	1526	1535	rVB	65886	104599	0.60%	0.040%
19	12.346	1552	1557	1561	rBV	72949	112508	0.65%	0.043%
20	12.416	1563	1569	1578	rVB	107664	188572	1.09%	0.072%
21	12.634	1600	1606	1611	rVB	61220	96829	0.56%	0.037%
22	12.857	1637	1644	1650	rBV3	72816	128702	0.74%	0.049%
23	13.110	1683	1687	1696	rVV3	39071	65871	0.38%	0.025%
24	13.404	1728	1737	1747	rVV2	127996	252574	1.45%	0.096%
25	13.769	1787	1799	1806	rBV7	40509	125831	0.72%	0.048%
26	13.910	1818	1823	1829	rVV	66350	113310	0.65%	0.043%
27	13.999	1829	1838	1845	rVV	6920967	9709550	55.92%	3.704%
28	14.057	1846	1848	1852	rVB3	50914	61094	0.35%	0.023%
29	14.110	1852	1857	1861	rBV4	45956	76966	0.44%	0.029%
30	14.175	1865	1868	1873	rVV2	41814	56354	0.32%	0.021%
31	14.281	1880	1886	1898	rBV2	8120089	12661464	72.91%	4.830%
32	14.475	1915	1919	1926	rVB	86642	125064	0.72%	0.048%
33	14.587	1928	1938	1945	rBV	4815904	7255769	41.78%	2.768%
34	14.651	1945	1949	1958	rVB2	100408	181706	1.05%	0.069%
35	14.787	1965	1972	1986	rBV	2615841	3958273	22.79%	1.510%
36	14.987	2001	2006	2012	rVB	194411	281126	1.62%	0.107%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	15.257	2048	2052	2056	rVB2	37247	51491	0.30%	0.020%
38	15.375	2066	2072	2075	rBV5	47582	82690	0.48%	0.032%
39	15.422	2076	2080	2084	rVB3	47529	68400	0.39%	0.026%
40	15.581	2100	2107	2113	rBV	9756568	14512139	83.57%	5.536%
41	15.634	2113	2116	2121	rVB	146530	215977	1.24%	0.082%
42	15.704	2122	2128	2136	rBV	2747419	3722491	21.44%	1.420%
43	15.828	2142	2149	2156	rBV6	56161	115537	0.67%	0.044%
44	15.951	2163	2170	2181	rVB	4978972	6883332	39.64%	2.626%
45	16.081	2188	2192	2198	rVB2	58384	102571	0.59%	0.039%
46	16.292	2223	2228	2230	rBV2	127619	190594	1.10%	0.073%
47	16.434	2248	2252	2261	rBV5	107505	206482	1.19%	0.079%
48	16.516	2261	2266	2274	rVB	186217	272974	1.57%	0.104%
49	16.651	2285	2289	2293	rVB5	38923	63194	0.36%	0.024%
50	16.728	2299	2302	2310	rVB3	80483	122524	0.71%	0.047%
51	16.869	2323	2326	2331	rBV3	43133	68042	0.39%	0.026%
52	16.998	2343	2348	2354	rBV2	131102	219972	1.27%	0.084%
53	17.087	2357	2363	2367	rVV4	82485	146388	0.84%	0.056%
54	17.163	2373	2376	2382	rVB	99892	144989	0.83%	0.055%
55	17.228	2383	2387	2394	rBV	212860	308986	1.78%	0.118%
56	17.292	2394	2398	2401	rBV2	161106	200265	1.15%	0.076%
57	17.345	2401	2407	2411	rVV	5583665	7687363	44.27%	2.932%
58	17.387	2411	2414	2419	rVV	1820194	2517204	14.50%	0.960%
59	17.445	2419	2424	2436	rVB	9783113	14438187	83.15%	5.507%
60	17.622	2450	2454	2463	rBV	45660	120803	0.70%	0.046%
61	17.751	2473	2476	2479	rVV	52899	60702	0.35%	0.023%
62	17.828	2486	2489	2492	rBV2	54964	62375	0.36%	0.024%
63	17.863	2492	2495	2500	rVB3	88385	104562	0.60%	0.040%
64	17.922	2501	2505	2509	rBV3	49270	86182	0.50%	0.033%
65	18.104	2532	2536	2541	rBV	159602	222368	1.28%	0.085%
66	18.163	2541	2546	2550	rBV	532968	735824	4.24%	0.281%
67	18.222	2550	2556	2560	rBV2	3186593	4302694	24.78%	1.641%
68	18.334	2573	2575	2581	rVB	138374	157591	0.91%	0.060%
69	18.398	2581	2586	2590	rBV3	338740	588972	3.39%	0.225%
70	18.504	2599	2604	2608	rBV3	223808	350376	2.02%	0.134%
71	18.692	2632	2636	2640	rVV	203528	268783	1.55%	0.103%
72	18.734	2640	2643	2651	rVB2	129191	190189	1.10%	0.073%
73	19.098	2700	2705	2710	rBV2	211847	424321	2.44%	0.162%
74	19.351	2738	2748	2755	rBV	4685960	6641754	38.25%	2.533%
75	19.445	2761	2764	2769	rBV	484628	600050	3.46%	0.229%
76	19.498	2769	2773	2777	rVB	349829	434294	2.50%	0.166%

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 Title : SVOA CALIBRATION

77	19.586	2785	2788	2796	rVB3	153550	275100	1.58%	0.105%
78	19.681	2798	2804	2807	rVV	10503450	15390512	88.63%	5.871%
79	19.704	2807	2808	2815	rVV	3448163	3137700	18.07%	1.197%
80	19.769	2816	2819	2825	rVB2	163217	228554	1.32%	0.087%
81	19.881	2834	2838	2843	rVB	178643	244408	1.41%	0.093%
82	20.016	2856	2861	2868	rBV3	247562	623532	3.59%	0.238%
83	20.186	2886	2890	2895	rVB2	702447	1021900	5.88%	0.390%
84	20.281	2903	2906	2910	rBV	352972	400495	2.31%	0.153%
85	20.475	2936	2939	2943	rBV	221706	330265	1.90%	0.126%
86	20.916	3011	3014	3019	rBV	362190	466953	2.69%	0.178%
87	21.080	3039	3042	3045	rBV	434660	477682	2.75%	0.182%
88	21.122	3045	3049	3053	rBV2	2000571	2890633	16.65%	1.103%
89	21.328	3080	3084	3088	rVB2	1090002	1307543	7.53%	0.499%
90	21.433	3095	3102	3106	rBV2	5469302	10841326	62.43%	4.135%
91	21.469	3106	3108	3118	rVB	2191356	2996429	17.26%	1.143%
92	23.151	3385	3394	3400	rBV2	2144993	7329794	42.21%	2.796%
93	23.692	3481	3486	3490	rBV	1204848	2377589	13.69%	0.907%
94	23.751	3490	3496	3502	rVV2	6570371	13894615	80.02%	5.300%
95	23.804	3502	3505	3512	rVB	1880260	3111304	17.92%	1.187%
96	23.910	3517	3523	3529	rBV2	3070854	6266046	36.08%	2.390%
97	24.421	3602	3610	3621	rVV2	7697606	17364791	100.00%	6.624%
98	24.527	3622	3628	3638	rVB	1865321	4238184	24.41%	1.617%
99	24.710	3655	3659	3674	rVB6	1062499	2470454	14.23%	0.942%
100	26.433	3947	3952	3959	rBV	1164217	3156726	18.18%	1.204%

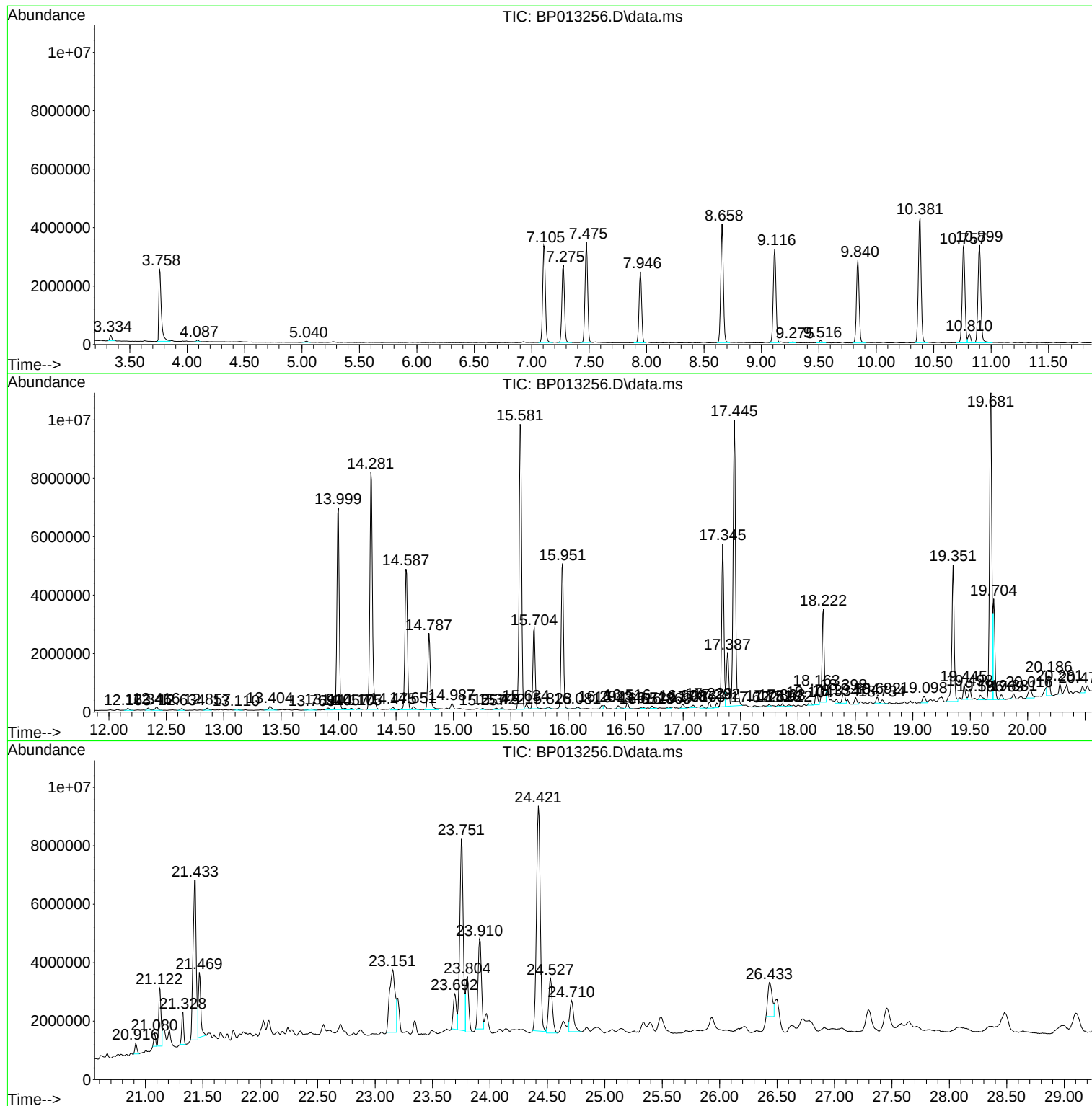
Sum of corrected areas: 262159347

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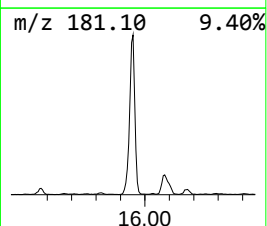
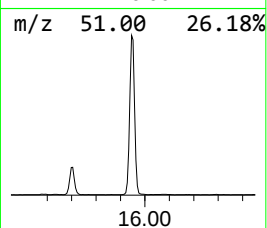
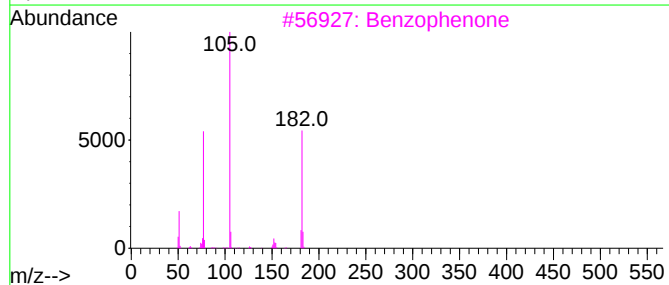
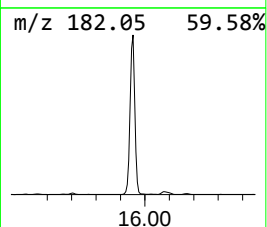
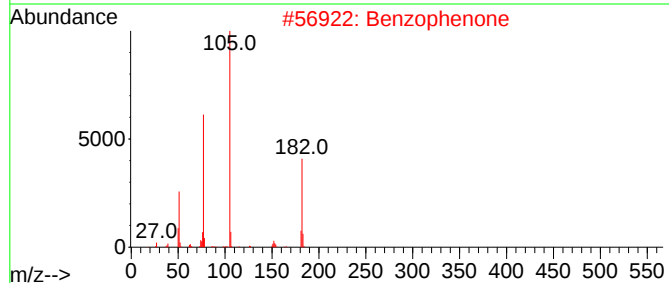
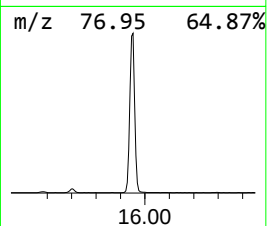
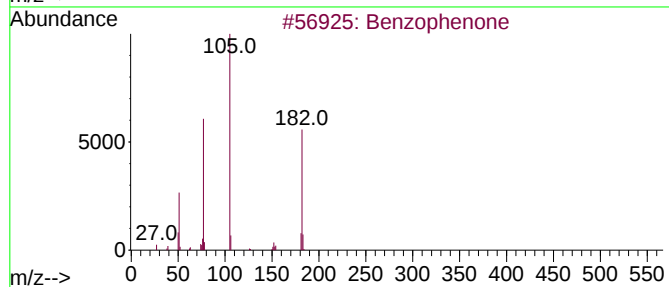
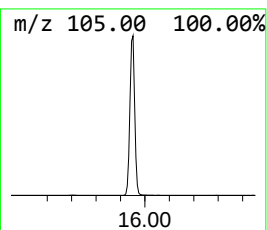
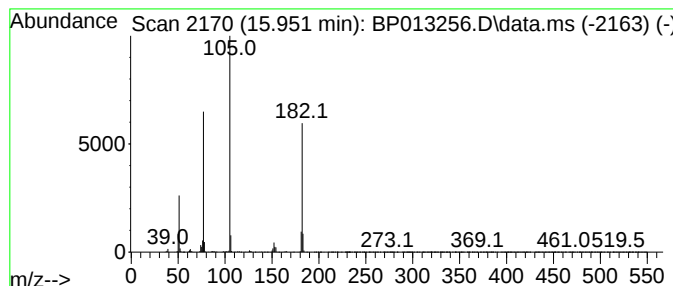
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.951	18.97 ng/ul	6883330	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	95
4			Benzophenone	182	C13H10O	000119-61-9	94
5			Benzenecarbothioic acid, S-methy...	152	C8H8OS	005925-68-8	49



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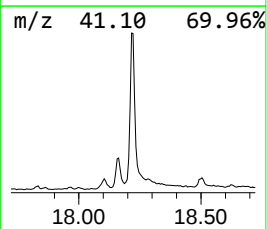
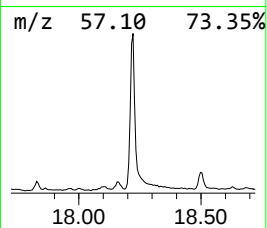
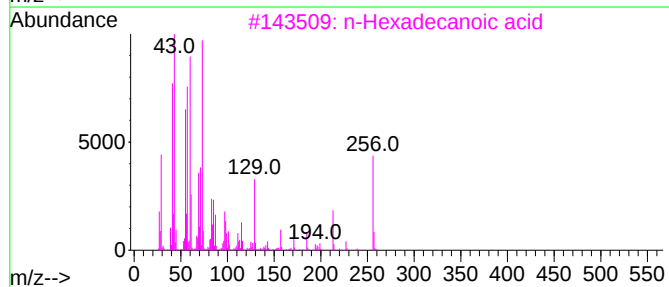
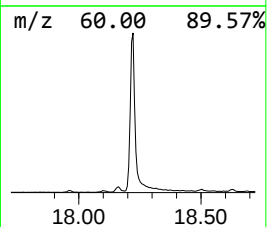
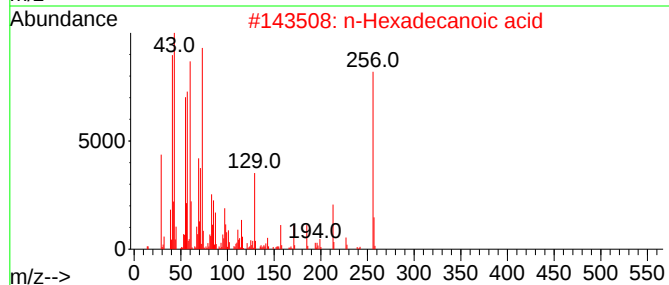
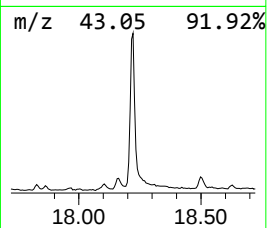
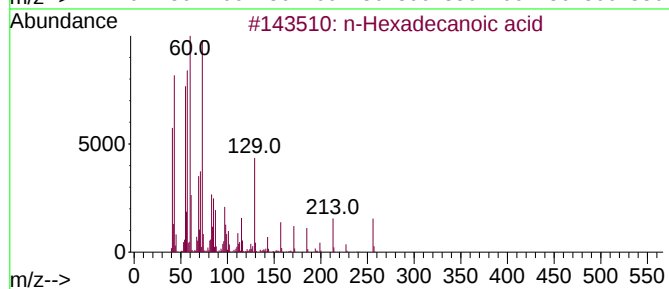
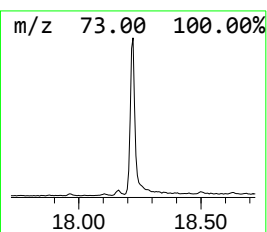
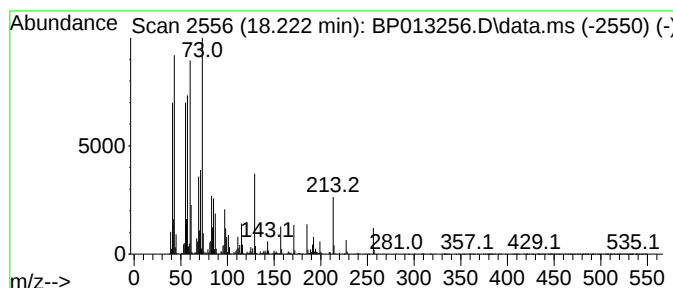
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.222	11.19 ng/ul	4302690	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	98
3			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
4			Tridecanoic acid	214	C13H26O2	000638-53-9	93
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	92



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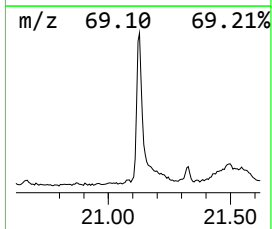
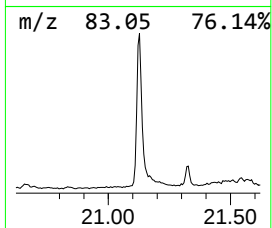
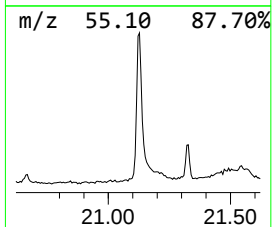
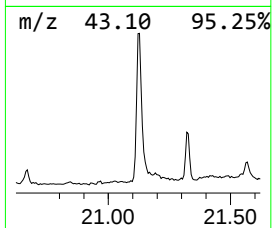
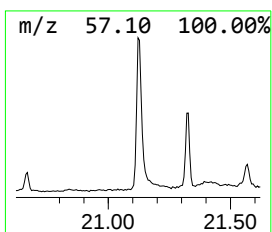
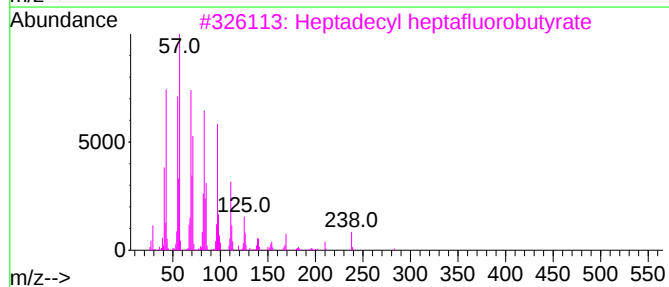
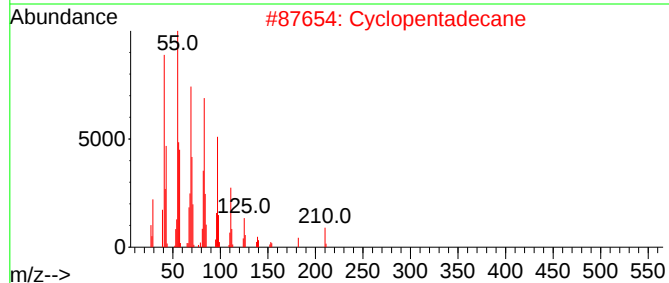
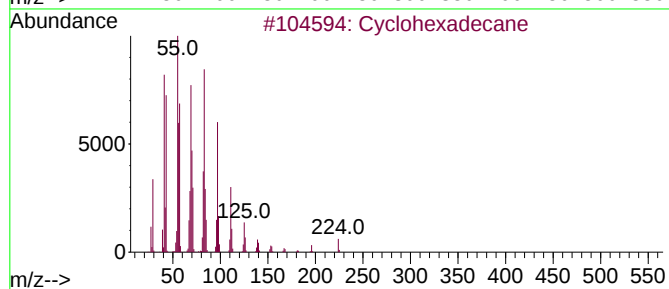
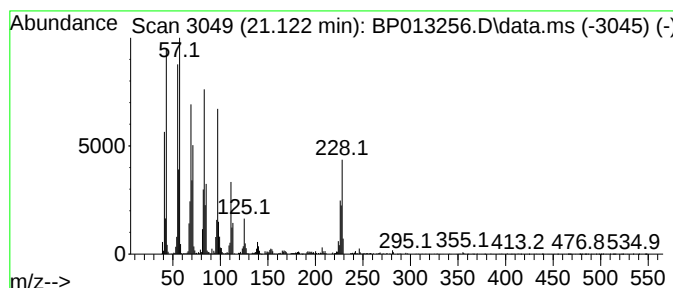
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 (DEL) Alkane: Cyclic21.122 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.122	5.33 ng/ul	2890630	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexadecane	224	C16H32	000295-65-8	92
2			Cyclopentadecane	210	C15H30	000295-48-7	90
3			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	89
4			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	89
5			1-Hexacosene	364	C26H52	018835-33-1	89



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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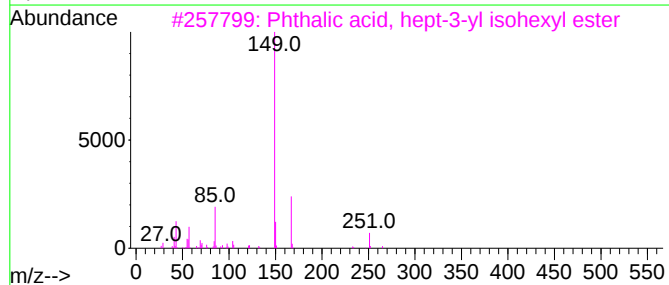
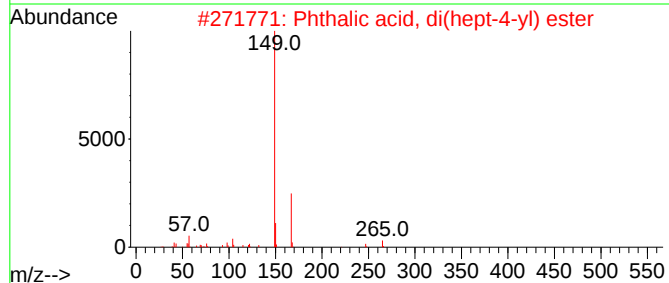
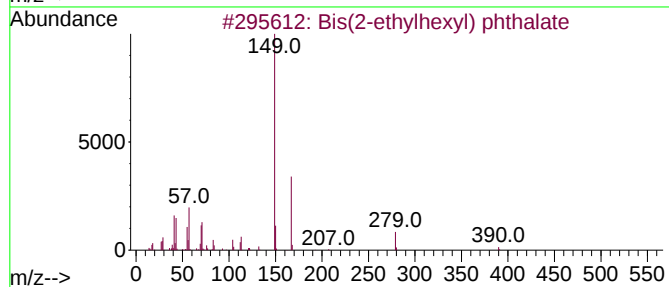
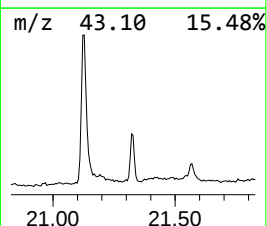
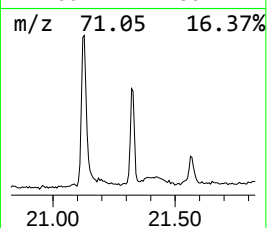
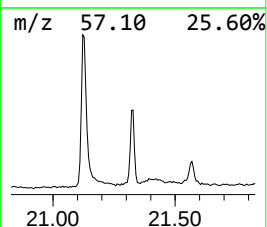
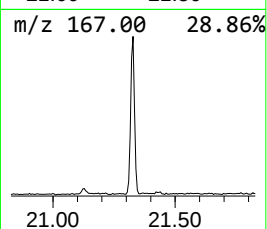
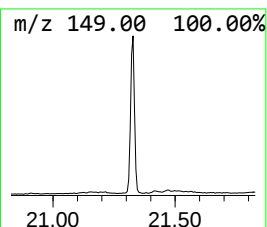
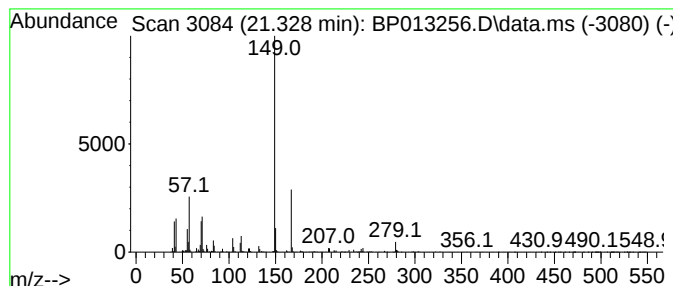
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Bis(2-ethylhexyl) phthalate Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.328	2.41 ng/ul	1307540	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bis(2-ethylhexyl) phthalate	390	C24H38O4	000117-81-7	90
2			Phthalic acid, di(hept-4-yl) ester	362	C22H34O4	1000356-80-0	64
3			Phthalic acid, hept-3-yl isohexy...	348	C21H32O4	1000356-98-9	64
4			1,2-Benzenedicarboxylic acid, bu...	334	C20H30O4	000085-69-8	59
5			Phthalic acid, cycloheptyl isoe...	346	C21H30O4	1000322-83-1	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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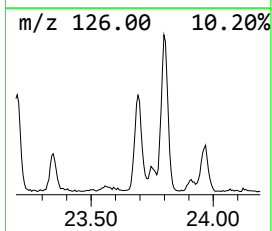
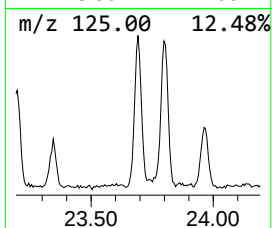
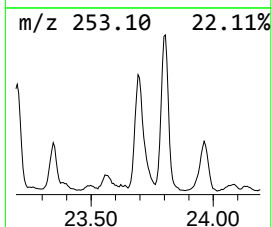
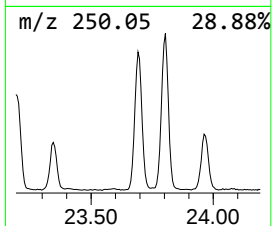
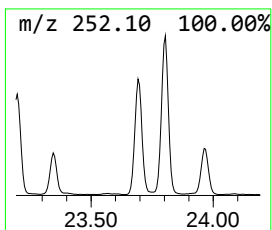
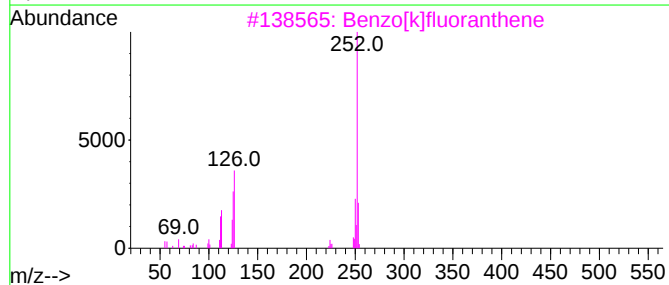
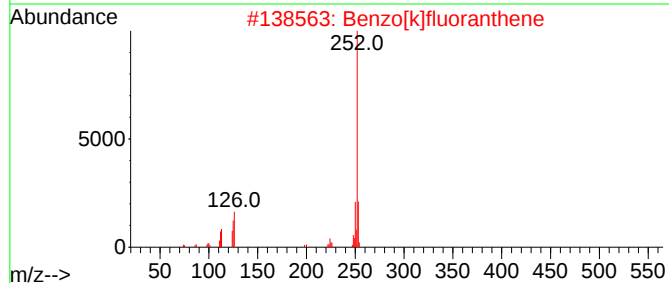
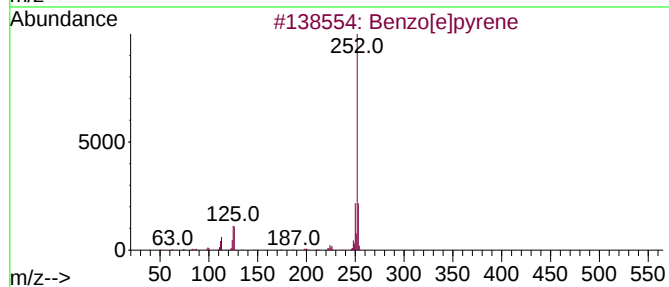
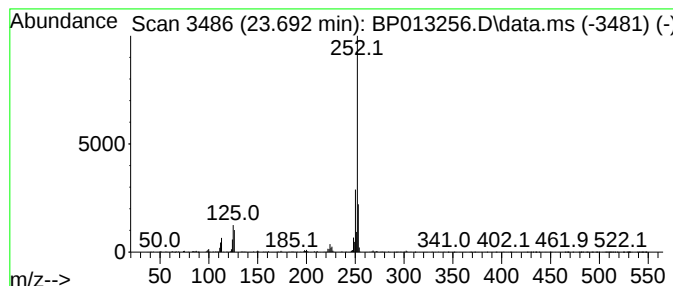
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzo[e]pyrene Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.692	7.59 ng/ul	2377590	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[e]pyrene	252	C20H12	000192-97-2	99
2			Benzo[k]fluoranthene	252	C20H12	000207-08-9	97
3			Benzo[k]fluoranthene	252	C20H12	000207-08-9	96
4			Benzo[b]fluoranthene	252	C20H12	000205-99-2	95
5			Benzo[a]pyrene	252	C20H12	000050-32-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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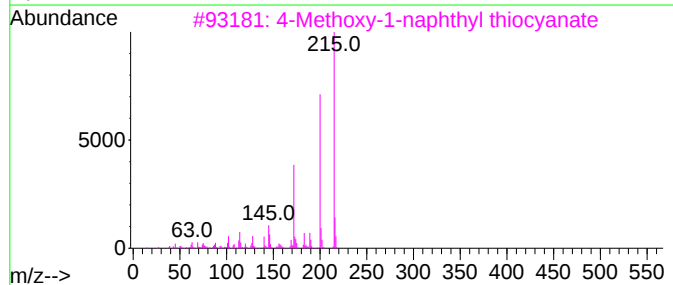
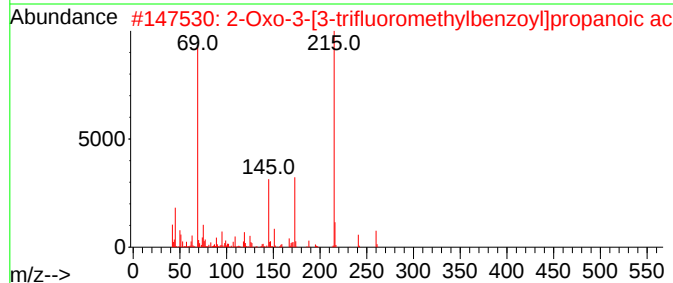
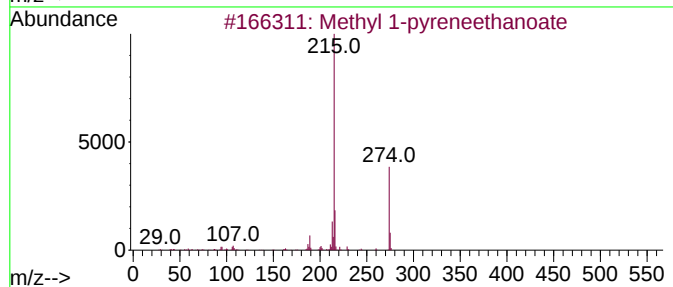
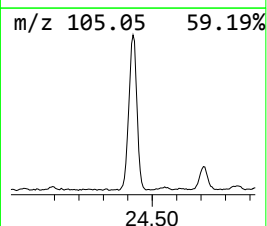
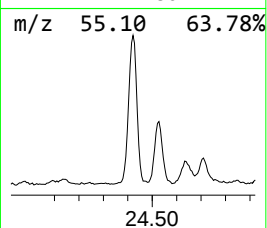
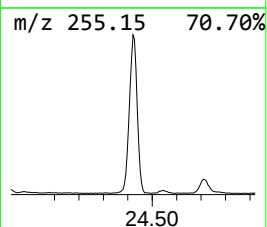
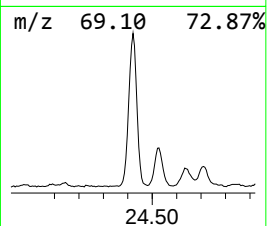
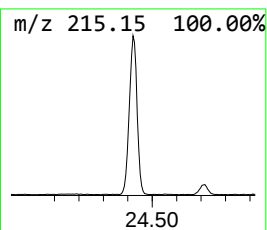
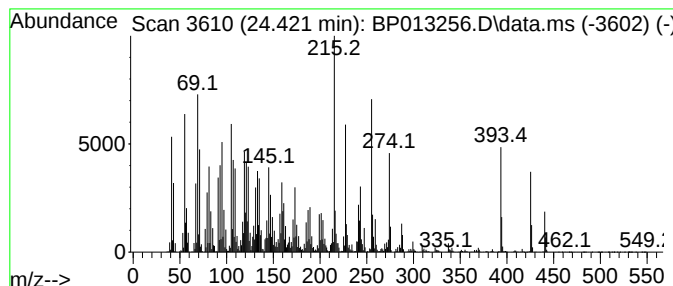
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.422	55.42 ng/ul	17364800	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	42
2			2-Oxo-3-[3-trifluoromethylbenzoyl]propanoic acid	260	C11H7F3O4	1000211-47-7	42
3			4-Methoxy-1-naphthyl thiocyanate	215	C12H9NOS	066231-77-4	20
4			1-Pyrenebutanoic acid	288	C20H16O2	003443-45-6	20
5			4-imidazolidinone, 1-methyl-2-(m...	215	C12H13N3O	1000396-19-4	18



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
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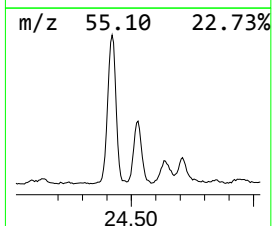
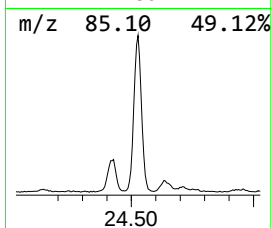
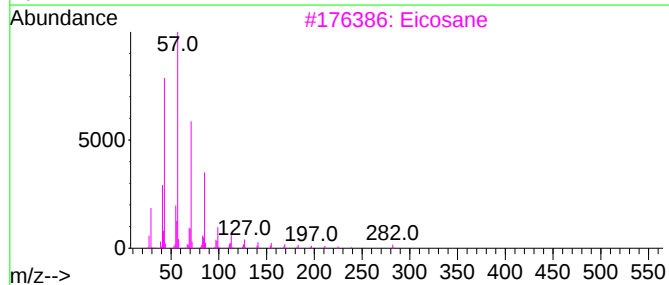
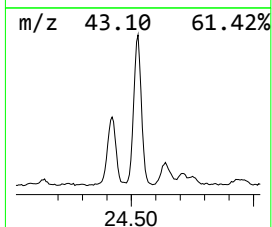
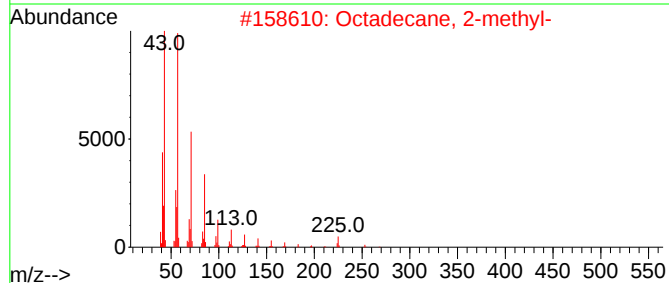
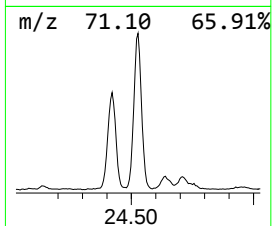
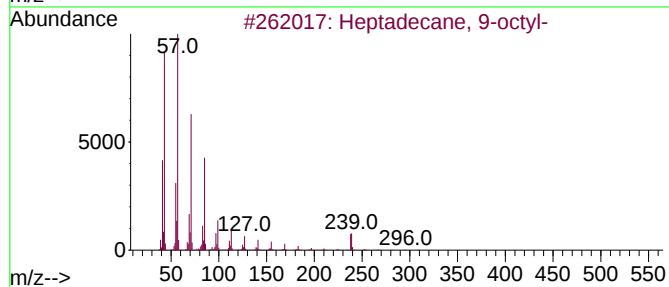
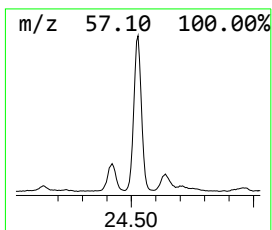
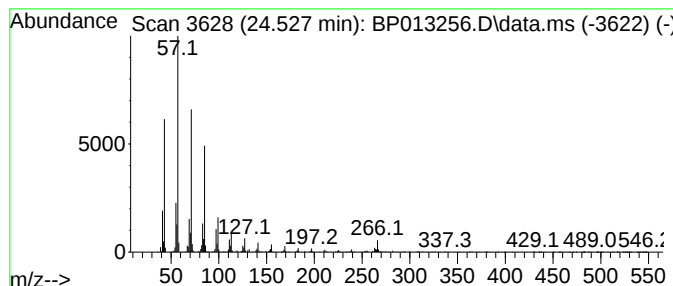
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.527	13.53 ng/ul	4238180	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	93
2	Octadecane, 2-methyl-		268	C19H40	001560-88-9	93
3	Eicosane		282	C20H42	000112-95-8	92
4	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	91
5	2-Bromo dodecane		248	C12H25Br	013187-99-0	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
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Sample : N6015-17
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Instrument :
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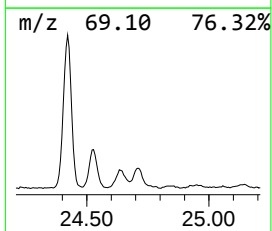
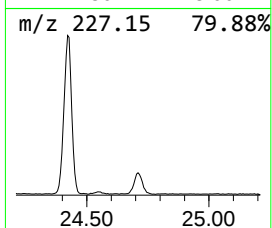
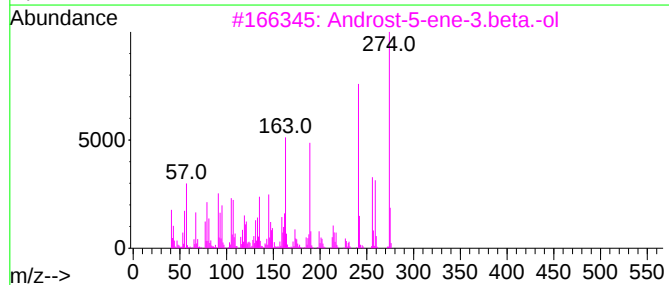
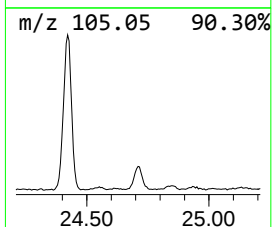
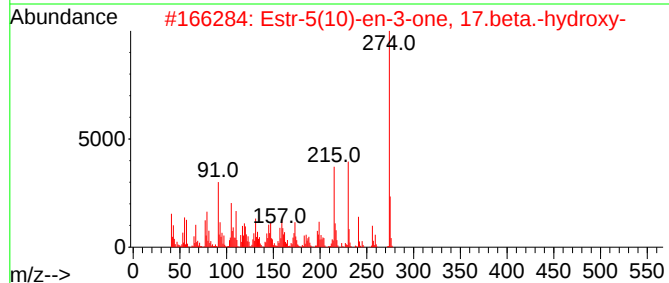
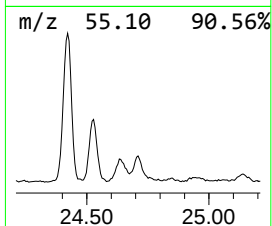
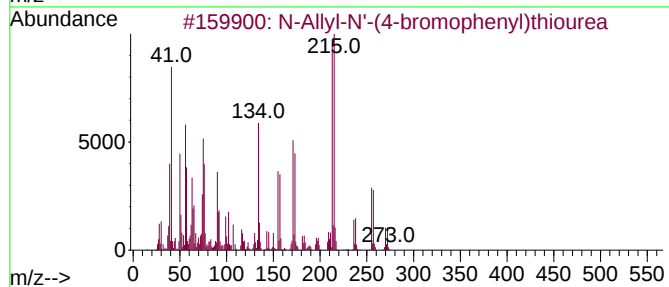
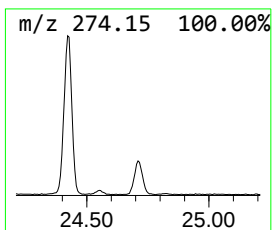
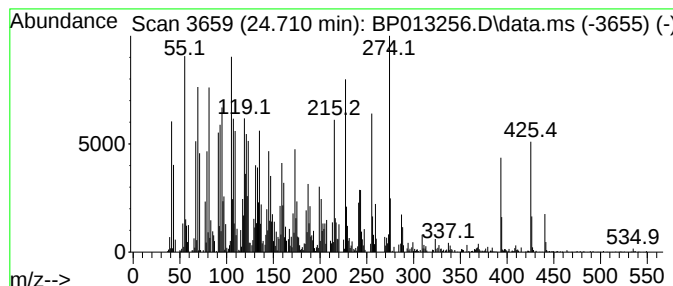
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 N-Allyl-N'-(4-bromophenyl)t... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.710	7.89 ng/ul	2470450	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Allyl-N'-(4-bromophenyl)thiourea	270	C10H11BrN2S	067544-10-9	90
2			Estr-5(10)-en-3-one, 17.beta.-hy...	274	C18H26O2	001089-78-7	42
3			Androst-5-ene-3.beta.-ol	274	C19H30O	001476-64-8	40
4			Methyl 1-pyreneethanoate	274	C19H14O2	073654-19-0	38
5			Androstan-15-one, (5.alpha.)-	274	C19H30O	000734-68-9	25



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
Misc :
ALS Vial : 23 Sample Multiplier: 1

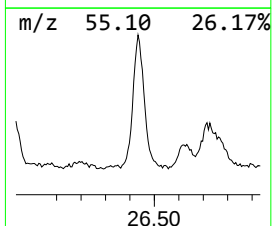
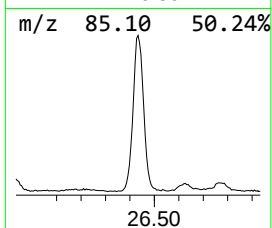
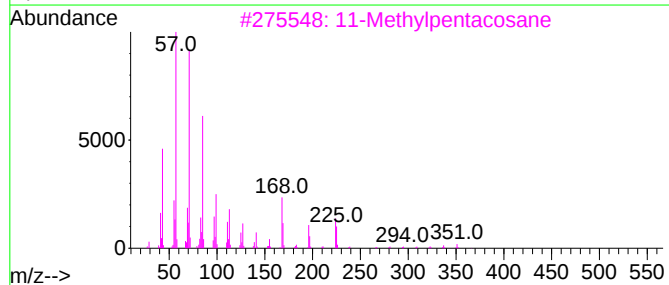
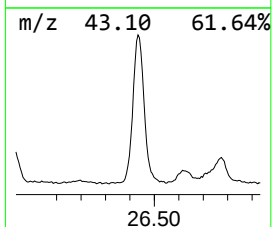
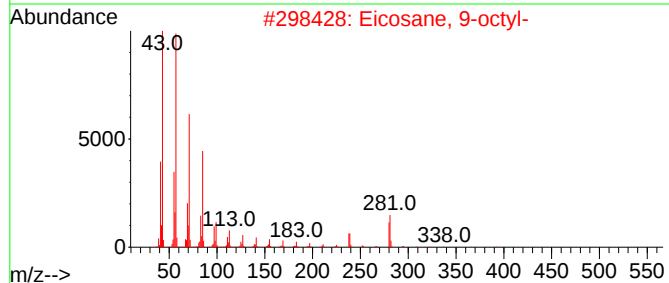
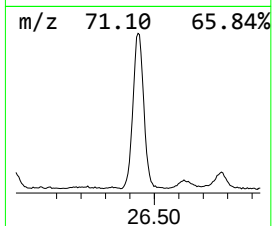
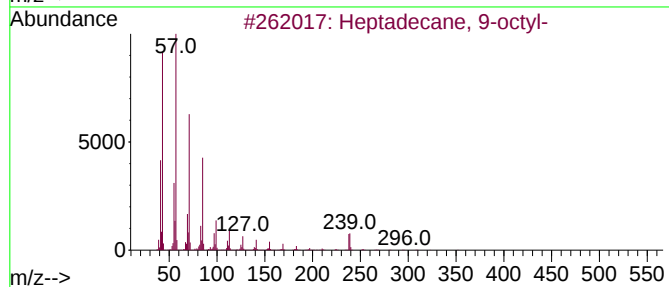
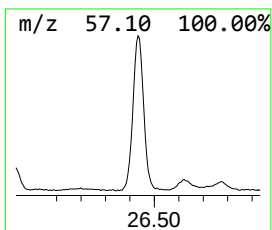
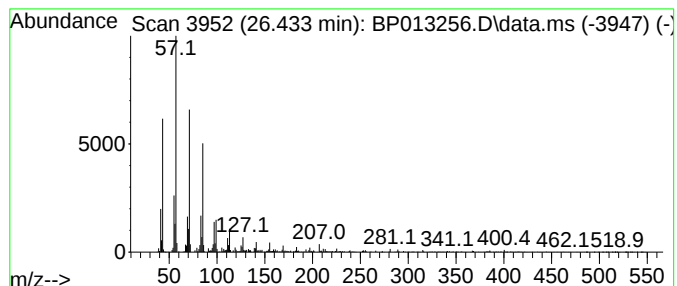
Instrument :
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD		R.T.	
26.433	10.08 ng/ul	3156730	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	94
2	Eicosane, 9-octyl-		394	C28H58	013475-77-9	93
3	11-Methylpentacosane		366	C26H54	015689-71-1	93
4	Heneicosane, 11-decyl-		437	C31H64	055320-06-4	90
5	Octacosane		394	C28H58	000630-02-4	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013256.D
Acq On : 22 Dec 2022 21:51
Operator : CG/JU
Sample : N6015-17
Misc :
ALS Vial : 23 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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzophenone	15.951	19.0	ng/u1	6883330	3	14.587	7255770	20.0
n-Hexadecanoic ...	18.222	11.2	ng/u1	4302690	4	17.345	7687360	20.0
(DEL) Alkane: C...	21.122	5.3	ng/u1	2890630	5	21.433	10841300	20.0
Bis(2-ethylhexy...	21.328	2.4	ng/u1	1307540	5	21.433	10841300	20.0
Benzo[e]pyrene	23.692	7.6	ng/u1	2377590	6	23.910	6266050	20.0
unknown-01	24.422	55.4	ng/u1	17364800	6	23.910	6266050	20.0
(DEL) Alkane: S...	24.527	13.5	ng/u1	4238180	6	23.910	6266050	20.0
N-Allyl-N'-(4-b...	24.710	7.9	ng/u1	2470450	6	23.910	6266050	20.0
(DEL) Alkane: S...	26.433	10.1	ng/u1	3156730	6	23.910	6266050	20.0