

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
 Data File : BP013247.D
 Acq On : 22 Dec 2022 15:43
 Operator : CG/JU
 Sample : N6015-06
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBXU3

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: BP013247.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	36	rVB	162787	228897	1.88%	0.145%
2	3.764	94	98	121	rBV	1504602	2253409	18.52%	1.427%
3	4.093	150	154	160	rVB	44081	61655	0.51%	0.039%
4	5.040	308	315	320	rVB2	36333	64973	0.53%	0.041%
5	5.264	348	353	359	rBV	35746	67995	0.56%	0.043%
6	6.546	567	571	577	rBV8	6827	13976	0.11%	0.009%
7	6.928	632	636	642	rVB8	14249	25192	0.21%	0.016%
8	7.105	659	666	679	rVB	2272216	3604099	29.62%	2.282%
9	7.275	688	695	705	rVB	1885074	2913198	23.94%	1.845%
10	7.375	709	712	718	rVB8	7726	13379	0.11%	0.008%
11	7.475	722	729	737	rBV	2392853	3721291	30.58%	2.356%
12	7.946	802	809	816	rBV	1932653	3007375	24.72%	1.904%
13	8.005	816	819	824	rVB6	13793	21406	0.18%	0.014%
14	8.446	890	894	898	rVB7	9675	13569	0.11%	0.009%
15	8.658	921	930	940	rBV	2553953	4113468	33.81%	2.605%
16	9.116	1000	1008	1020	rBV	2148507	3525788	28.98%	2.232%
17	9.269	1030	1034	1044	rVB7	21426	43178	0.35%	0.027%
18	9.516	1070	1076	1084	rBV2	55599	92794	0.76%	0.059%
19	9.681	1103	1104	1113	rVB8	9115	16315	0.13%	0.010%
20	9.840	1122	1131	1142	rBV	1774245	2960076	24.33%	1.874%
21	10.381	1215	1223	1234	rBV	2644308	4435277	36.45%	2.808%
22	10.534	1242	1249	1255	rVB10	13777	26472	0.22%	0.017%
23	10.763	1279	1288	1294	rBV	2407722	4027910	33.10%	2.550%
24	10.899	1304	1311	1327	rVV	1830718	3152974	25.91%	1.996%
25	11.175	1356	1358	1363	rVB6	9741	13382	0.11%	0.008%
26	11.287	1373	1377	1382	rVB8	13390	21492	0.18%	0.014%
27	11.540	1416	1420	1422	rBV4	9562	14746	0.12%	0.009%
28	12.016	1495	1501	1506	rBV7	12133	22634	0.19%	0.014%
29	12.163	1522	1526	1533	rVB	27256	41099	0.34%	0.026%
30	12.251	1537	1541	1547	rBV9	11086	18121	0.15%	0.011%
31	12.346	1552	1557	1561	rVB	41134	62192	0.51%	0.039%
32	12.746	1619	1625	1630	rBV9	9436	16249	0.13%	0.010%
33	12.798	1630	1634	1638	rVB7	9807	16811	0.14%	0.011%
34	12.910	1648	1653	1656	rBV7	9728	16479	0.14%	0.010%
35	13.116	1682	1688	1691	rVB7	15835	27253	0.22%	0.017%
36	13.328	1719	1724	1727	rBV7	10387	16405	0.13%	0.010%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	13.404	1730	1737	1742	rVB	53786	78791	0.65%	0.050%
38	13.510	1752	1755	1759	rBV5	22656	31401	0.26%	0.020%
39	13.793	1800	1803	1807	rVB6	13118	16218	0.13%	0.010%
40	13.857	1807	1814	1818	rBV10	17552	26630	0.22%	0.017%
41	13.922	1818	1825	1830	rBV	115143	170680	1.40%	0.108%
42	13.993	1830	1837	1845	rVV	3955419	5412040	44.48%	3.427%
43	14.057	1845	1848	1853	rVV7	33205	51041	0.42%	0.032%
44	14.110	1853	1857	1863	rVB9	23574	37877	0.31%	0.024%
45	14.175	1863	1868	1873	rVB2	19247	27757	0.23%	0.018%
46	14.228	1873	1877	1879	rBV4	11420	14782	0.12%	0.009%
47	14.281	1879	1886	1898	rVV	4692693	6908521	56.78%	4.374%
48	14.357	1898	1899	1903	rVB4	15810	14019	0.12%	0.009%
49	14.475	1916	1919	1926	rVB2	42596	79765	0.66%	0.051%
50	14.545	1926	1931	1932	rBV3	36564	44905	0.37%	0.028%
51	14.587	1932	1938	1947	rVB	3164280	4637370	38.11%	2.936%
52	14.787	1965	1972	1992	rBV	1357141	2245163	18.45%	1.422%
53	14.945	1995	1999	2004	rVV6	26378	43705	0.36%	0.028%
54	14.992	2004	2007	2014	rVV3	35603	51647	0.42%	0.033%
55	15.240	2044	2049	2056	rVB3	10671	22663	0.19%	0.014%
56	15.422	2078	2080	2085	rVB4	16841	23121	0.19%	0.015%
57	15.581	2101	2107	2114	rBV	5727659	8139638	66.89%	5.154%
58	15.698	2122	2127	2137	rBV	1467782	2023695	16.63%	1.281%
59	15.822	2144	2148	2155	rVB4	31627	62821	0.52%	0.040%
60	15.945	2164	2169	2181	rVB	1964730	2687755	22.09%	1.702%
61	16.145	2199	2203	2208	rVB8	16695	31514	0.26%	0.020%
62	16.210	2210	2214	2219	rBV6	33876	52306	0.43%	0.033%
63	16.292	2224	2228	2236	rBV8	23021	49399	0.41%	0.031%
64	16.434	2248	2252	2254	rBV5	24462	33470	0.28%	0.021%
65	16.516	2261	2266	2271	rBV2	73891	104280	0.86%	0.066%
66	16.728	2298	2302	2306	rBV3	53352	77293	0.64%	0.049%
67	17.069	2355	2360	2362	rBV2	35710	61453	0.51%	0.039%
68	17.228	2383	2387	2391	rBV3	65595	96308	0.79%	0.061%
69	17.292	2395	2398	2401	rBV2	41993	52309	0.43%	0.033%
70	17.345	2401	2407	2411	rBV	3605504	5159217	42.40%	3.267%
71	17.445	2417	2424	2432	rVB	5688561	8004190	65.78%	5.068%
72	17.557	2439	2443	2450	rVB3	69972	83495	0.69%	0.053%
73	17.698	2463	2467	2476	rBV3	145469	219648	1.81%	0.139%
74	18.004	2516	2519	2521	rBV3	32975	36659	0.30%	0.023%
75	18.104	2532	2536	2539	rBV2	66458	87245	0.72%	0.055%
76	18.157	2542	2545	2550	rBV2	100386	139158	1.14%	0.088%

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77	18.216	2550	2555	2574	rVV	1275968	1992858	16.38%	1.262%
78	18.504	2601	2604	2608	rBV5	39647	54644	0.45%	0.035%
79	18.545	2608	2611	2616	rVV4	46693	67255	0.55%	0.043%
80	18.628	2622	2625	2628	rBV4	43540	60327	0.50%	0.038%
81	18.781	2648	2651	2656	rBV4	64810	84004	0.69%	0.053%
82	19.116	2703	2708	2712	rBV3	80353	146987	1.21%	0.093%
83	19.204	2719	2723	2728	rBV	105426	162113	1.33%	0.103%
84	19.251	2728	2731	2735	rVB	92947	110703	0.91%	0.070%
85	19.333	2737	2745	2746	rBV3	189615	437280	3.59%	0.277%
86	19.445	2760	2764	2773	rBV	323767	427775	3.52%	0.271%
87	19.680	2798	2804	2812	rBV	7358115	10170955	83.59%	6.440%
88	19.828	2824	2829	2835	rBV	1645484	1935239	15.90%	1.225%
89	19.951	2847	2850	2855	rBV2	101198	185102	1.52%	0.117%
90	20.175	2884	2888	2894	rVB4	243913	454567	3.74%	0.288%
91	21.127	3045	3050	3053	rBV	2110172	3612052	29.69%	2.287%
92	21.433	3097	3102	3106	rBV	4832883	6857379	56.36%	4.342%
93	22.039	3202	3205	3207	rBV	331443	430150	3.54%	0.272%
94	23.121	3385	3389	3395	rBV	1005801	1590791	13.07%	1.007%
95	23.751	3489	3496	3511	rVB2	5843497	12167904	100.00%	7.704%
96	23.910	3517	3523	3533	rVB2	3196308	6474869	53.21%	4.100%
97	24.521	3623	3627	3638	rVB	1858556	4048954	33.28%	2.564%
98	25.933	3862	3867	3879	rVB	1081269	2689289	22.10%	1.703%
99	26.439	3945	3953	3971	rVB	2618113	8529938	70.10%	5.401%
100	29.004	4380	4389	4401	rBV6	2571629	9409928	77.33%	5.958%

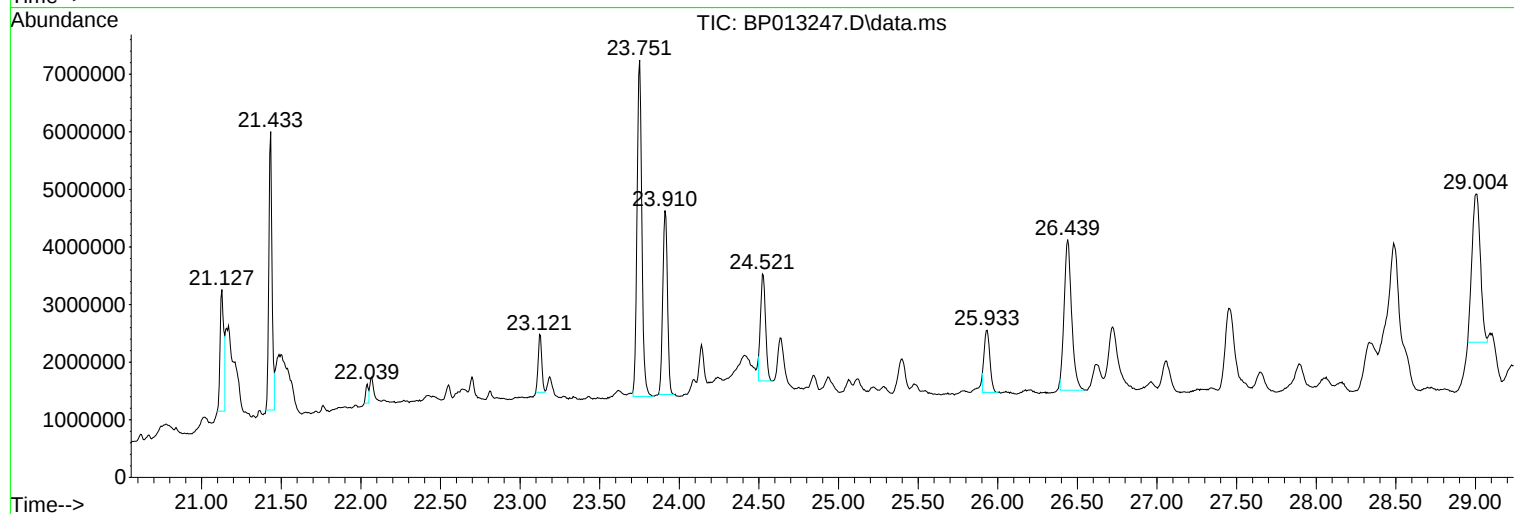
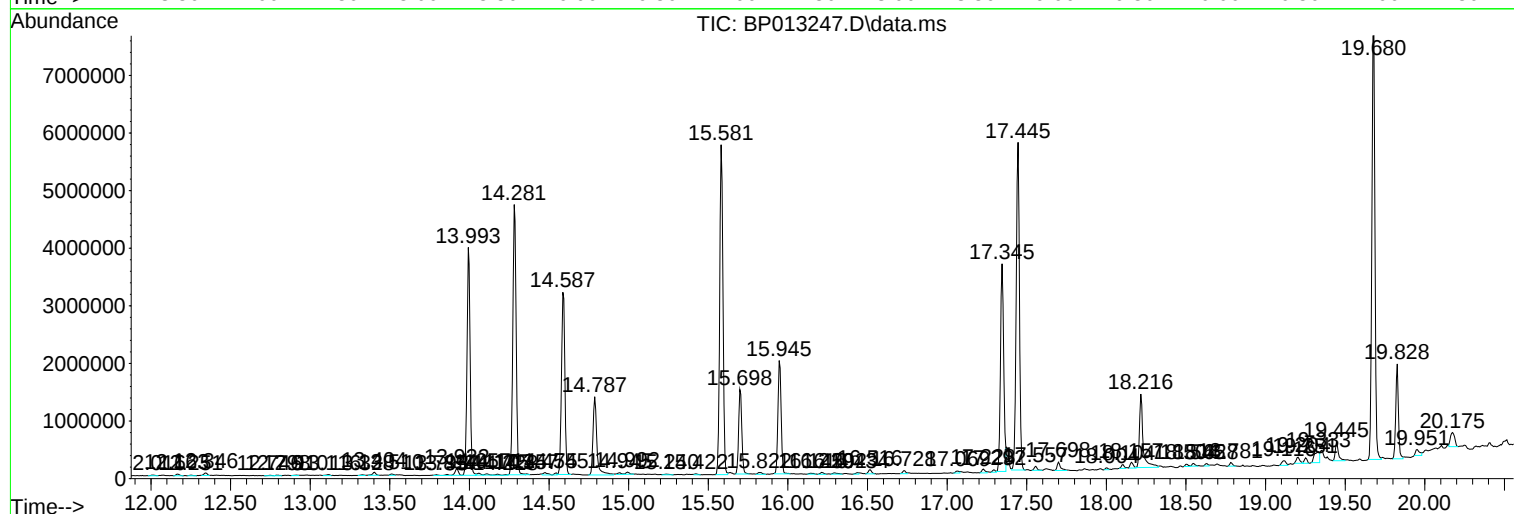
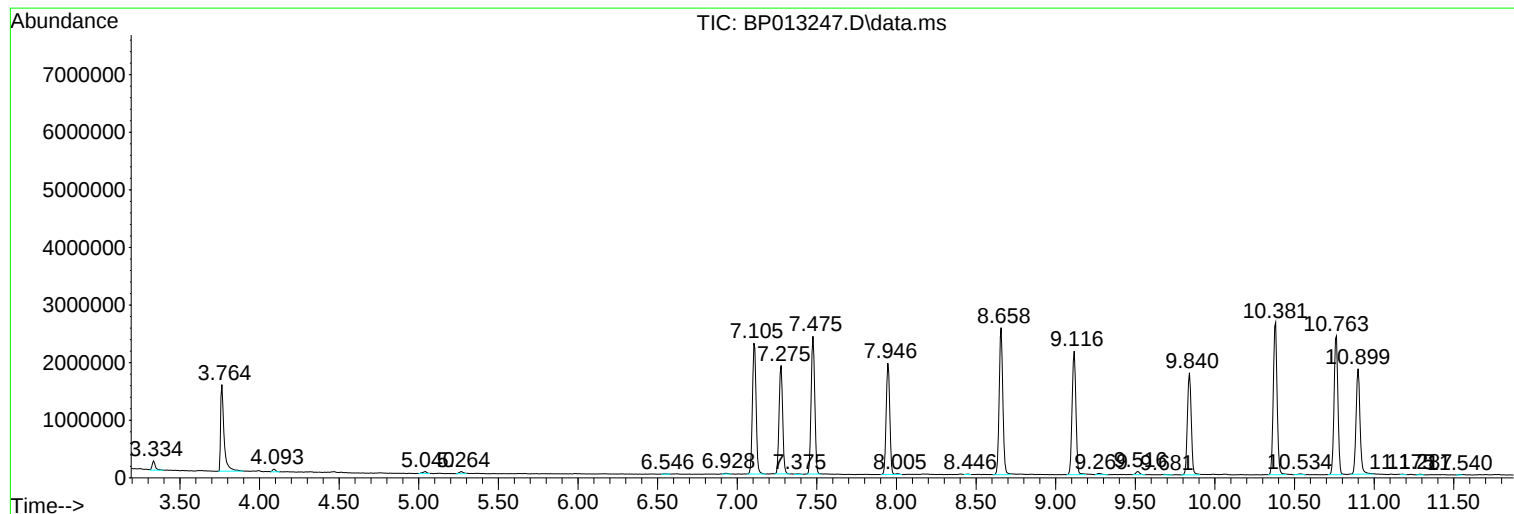
Sum of corrected areas: 157932541

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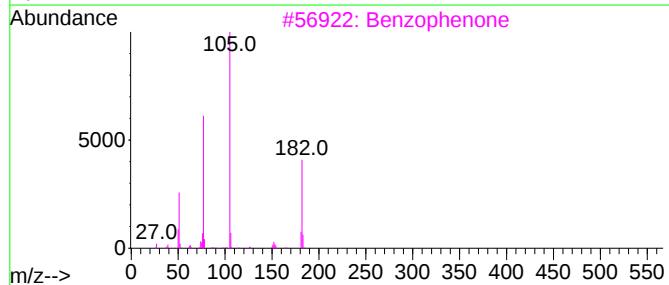
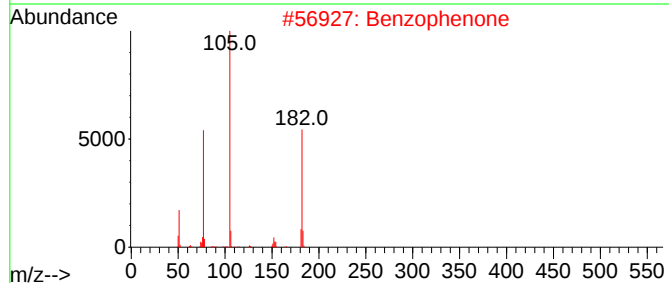
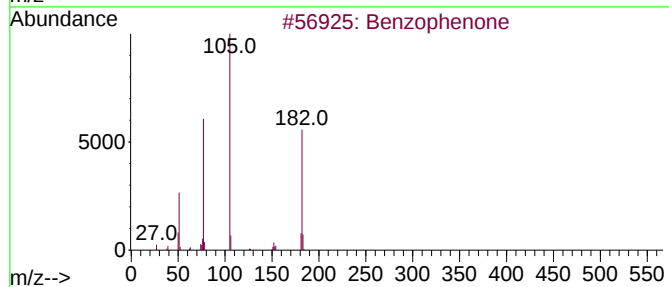
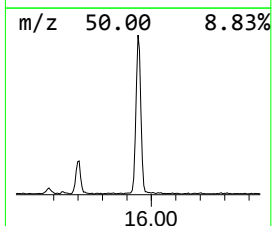
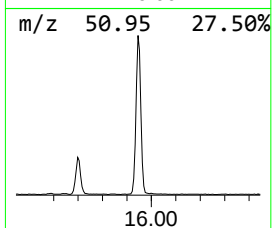
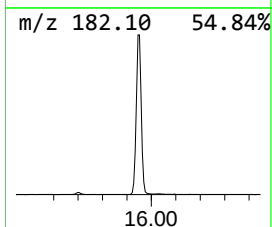
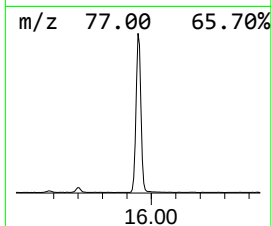
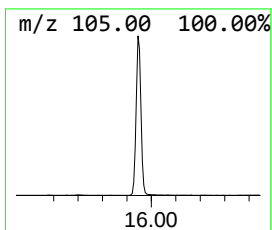
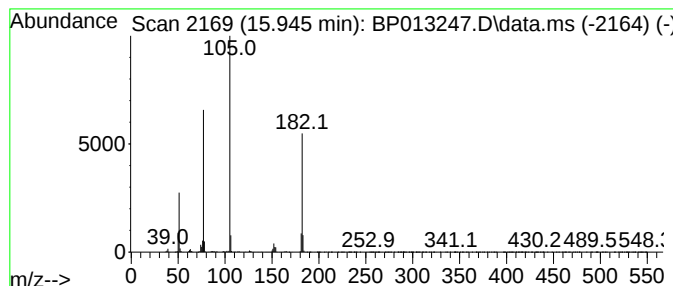
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	11.59 ng/ul	2687760	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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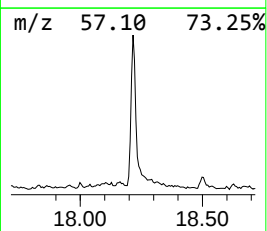
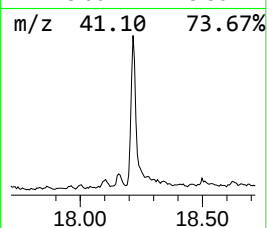
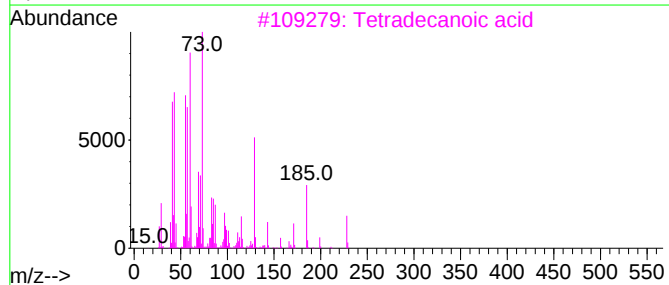
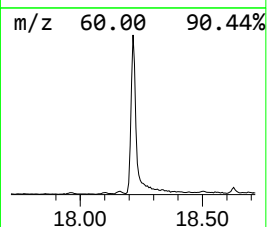
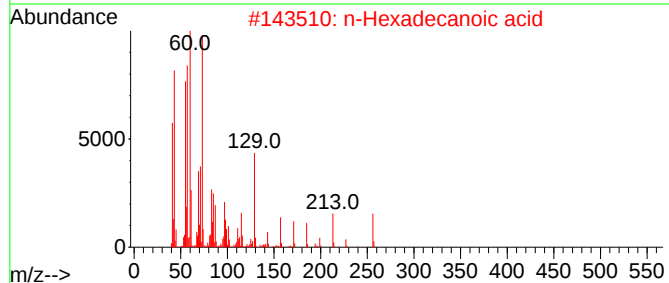
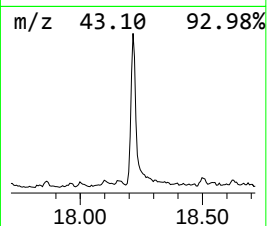
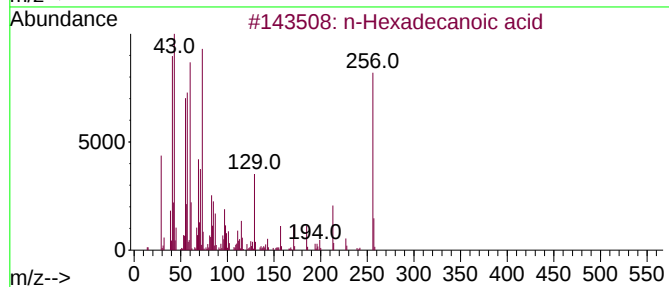
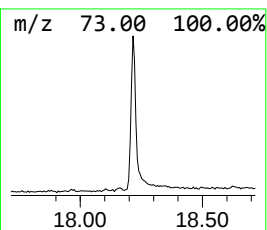
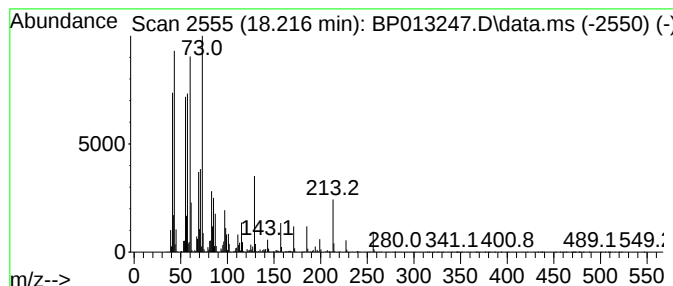
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	7.73 ng/ul	1992860	Phenanthrene-d10	17.345

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
2		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	98
3		Tetradecanoic acid	228	C14H28O2	000544-63-8	93
4		n-Hexadecanoic acid	256	C16H32O2	000057-10-3	87
5		Pentadecanoic acid	242	C15H30O2	001002-84-2	87



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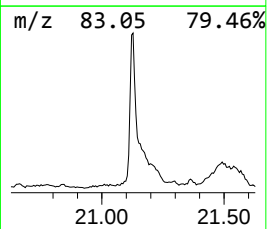
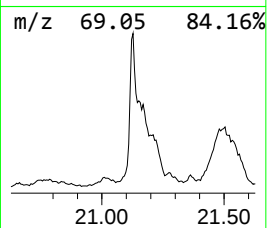
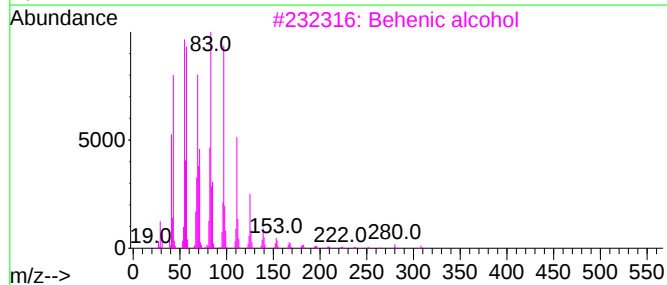
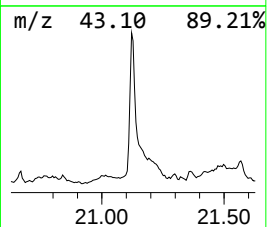
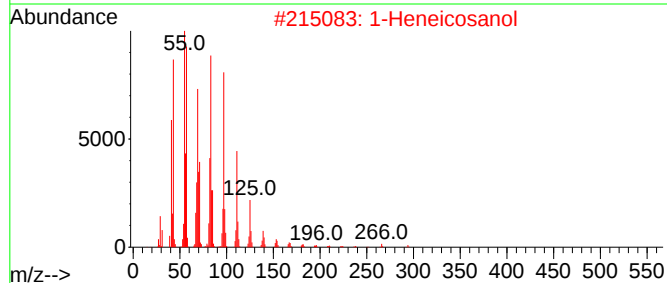
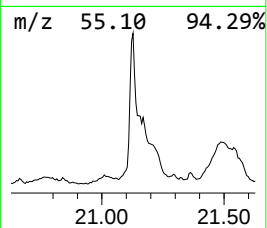
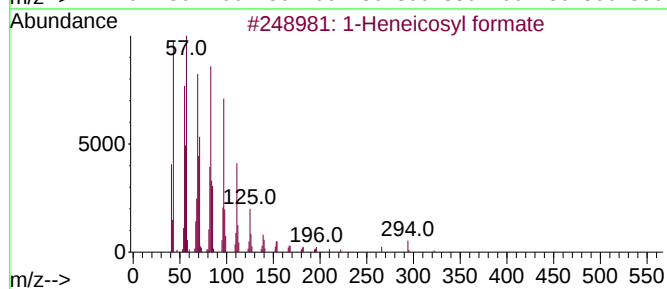
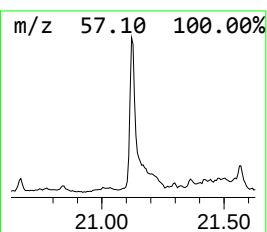
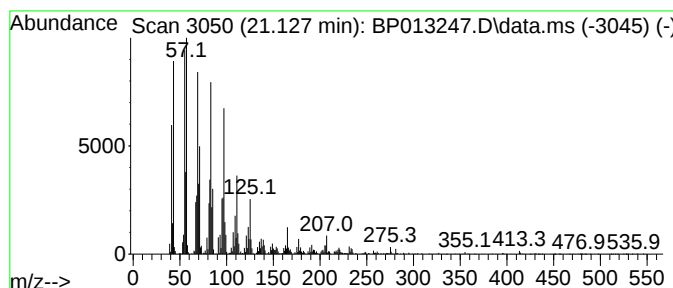
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-Heneicosyl formate Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.	
21.128	10.53 ng/ul	3612050	Chrysene-d12	21.433	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Heneicosyl formate	340	C22H44O2	077899-03-7	97
2	1-Heneicosanol	312	C21H44O	015594-90-8	93
3	Behenic alcohol	326	C22H46O	000661-19-8	91
4	Octacosanol	410	C28H58O	000557-61-9	90
5	Heptacos-1-ene	378	C27H54	015306-27-1	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
Operator : CG/JU
Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXU3

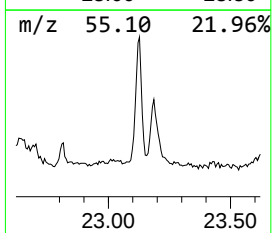
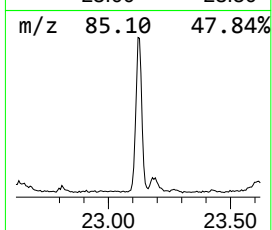
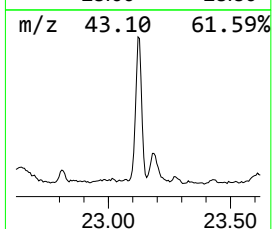
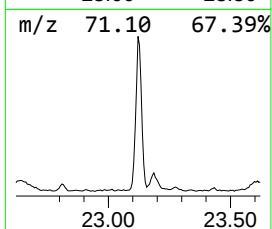
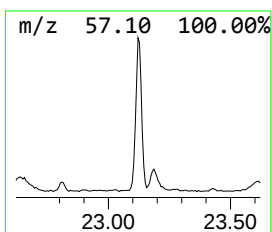
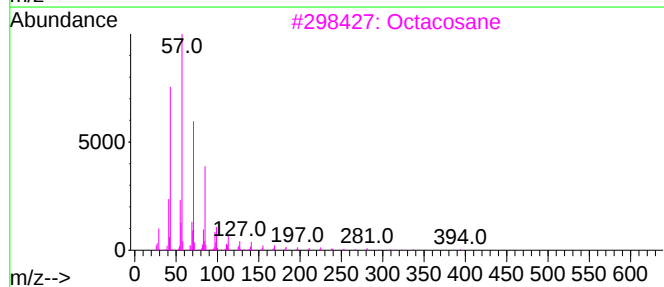
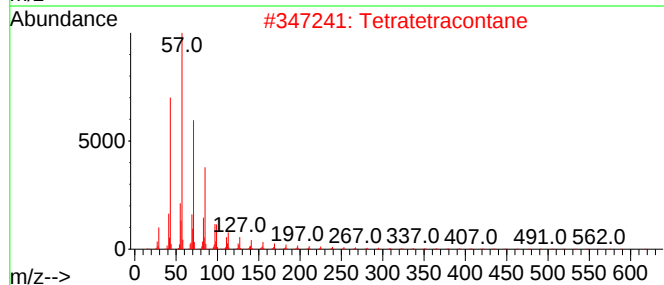
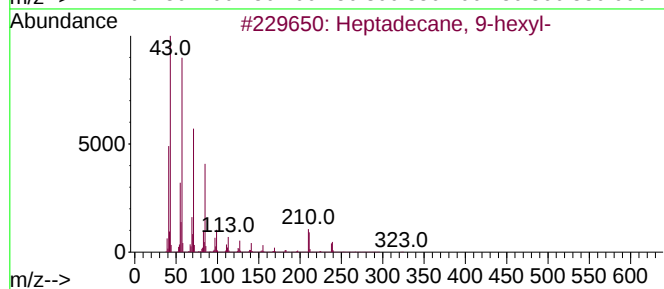
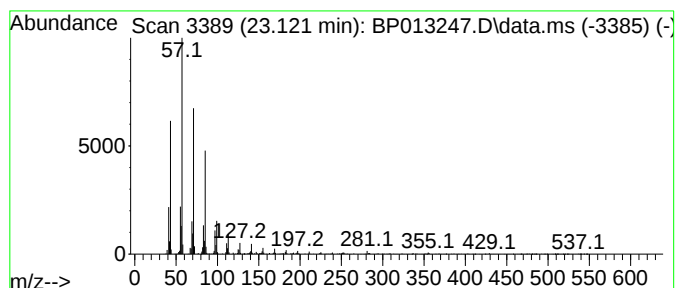
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.122	4.91 ng/ul	1590790	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Heptadecane, 9-hexyl-	324	C23H48	055124-79-3	94
2			Tetratetracontane	619	C44H90	007098-22-8	91
3			Octacosane	394	C28H58	000630-02-4	91
4			Octacosane, 2-methyl-	408	C29H60	001560-98-1	91
5			Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
Operator : CG/JU
Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

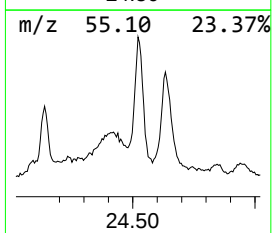
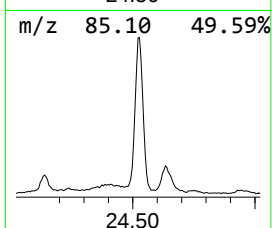
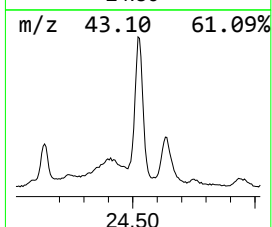
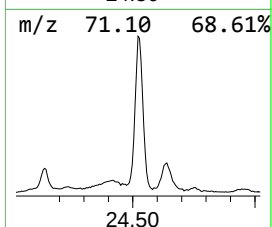
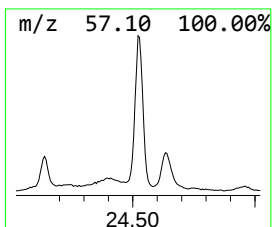
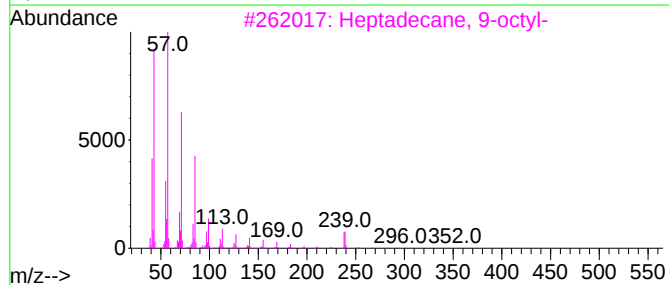
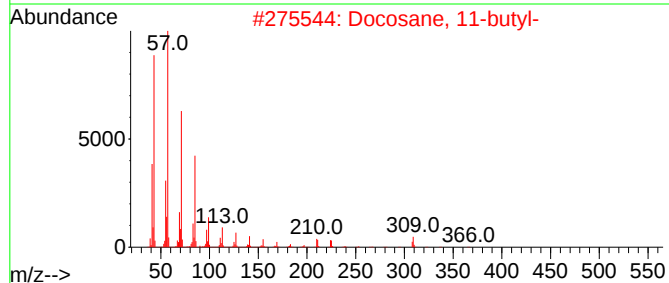
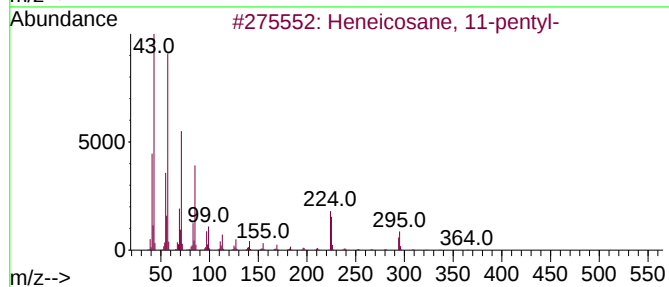
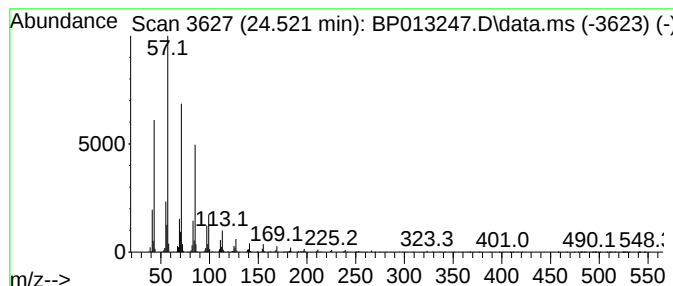
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.521	12.51 ng/ul	4048950	Perylene-d12		23.910	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Heneicosane, 11-pentyl-		366	C26H54	014739-72-1	94
2	Docosane, 11-butyl-		366	C26H54	013475-76-8	94
3	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91
4	Dotriacontane, 1-iodo-		576	C32H65I	1000406-32-4	91
5	Docosane, 1-iodo-		436	C22H45I	1000406-31-9	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
Operator : CG/JU
Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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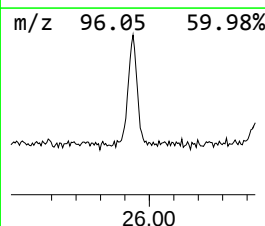
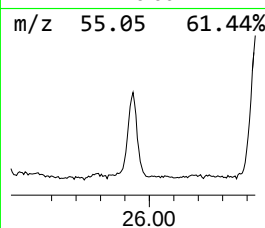
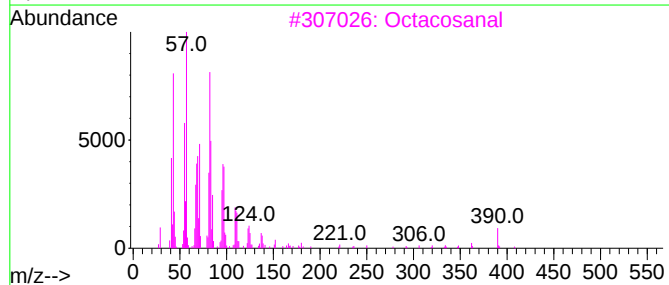
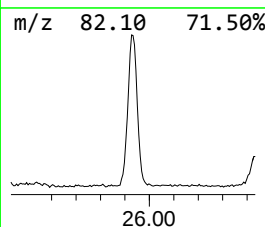
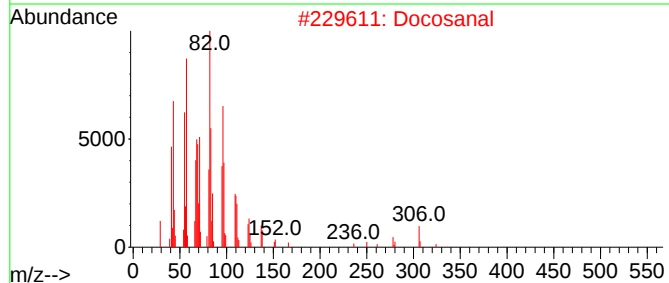
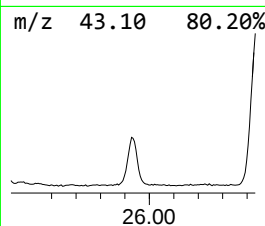
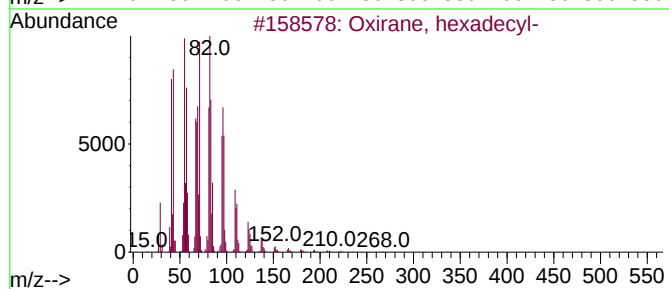
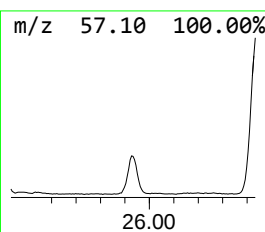
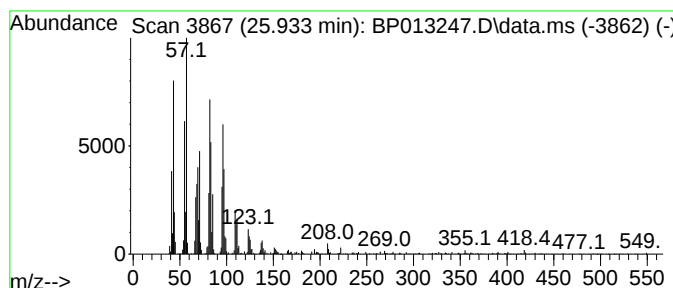
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Oxirane, hexadecyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
25.933	8.31 ng/ul	2689290	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	91
2			Docosanal	324	C22H44O	057402-36-5	87
3			Octacosanal	408	C28H56O	022725-64-0	83
4			(Z)-14-Tricosenyl formate	366	C24H46O2	077899-10-6	70
5			Oxirane, hexadecyl-	268	C18H36O	007390-81-0	68



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
Operator : CG/JU
Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXU3

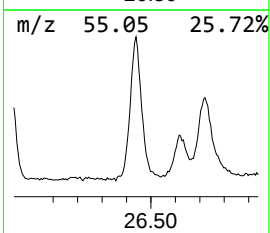
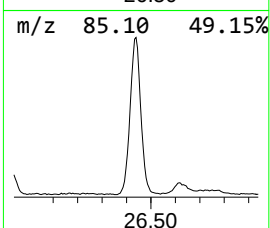
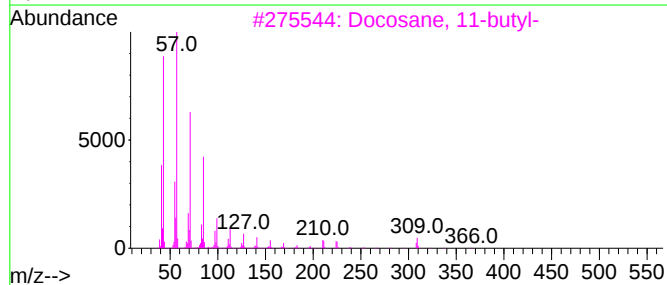
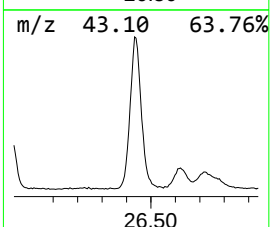
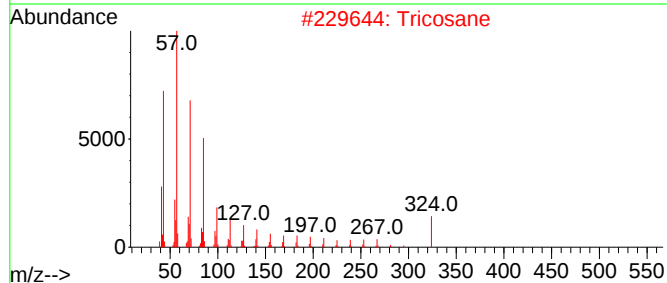
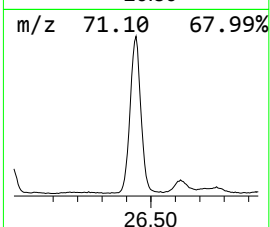
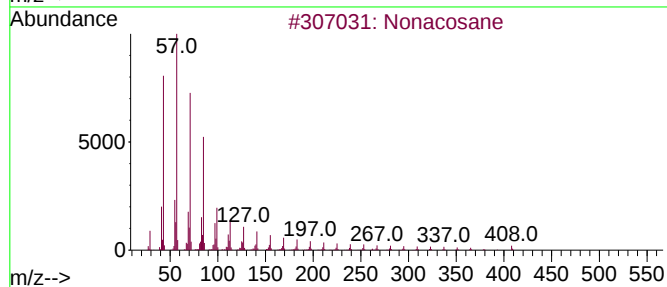
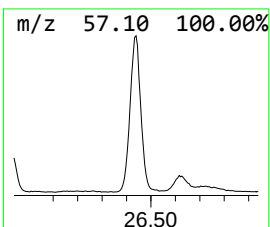
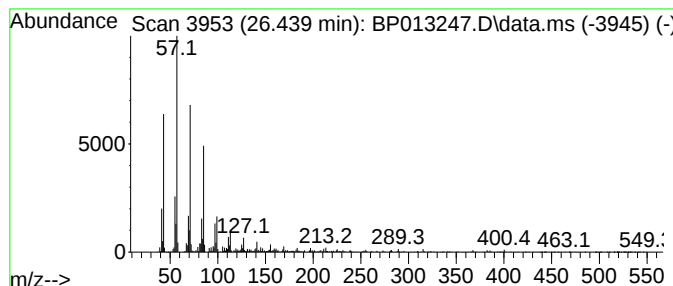
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 (DEL) Alkane: Straight-Chai... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
26.439	26.35 ng/ul	8529940	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Nonacosane	408	C29H60	000630-03-5	95
2			Tricosane	324	C23H48	000638-67-5	95
3			Docosane, 11-butyl-	366	C26H54	013475-76-8	95
4			Docosane, 7-hexyl-	394	C28H58	055373-86-9	93
5			Eicosane, 10-methyl-	296	C21H44	054833-23-7	93



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
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Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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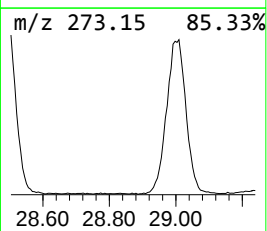
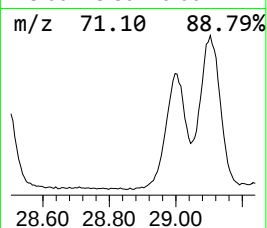
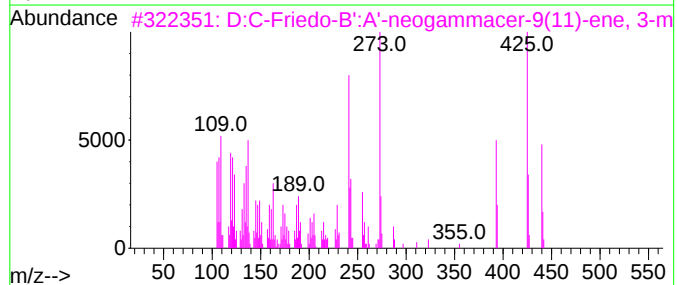
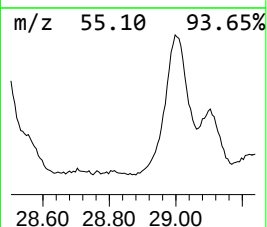
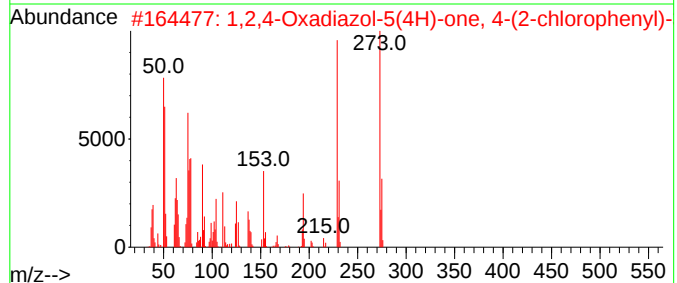
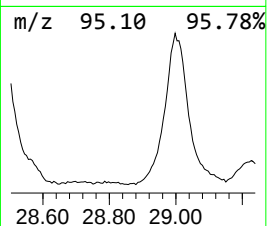
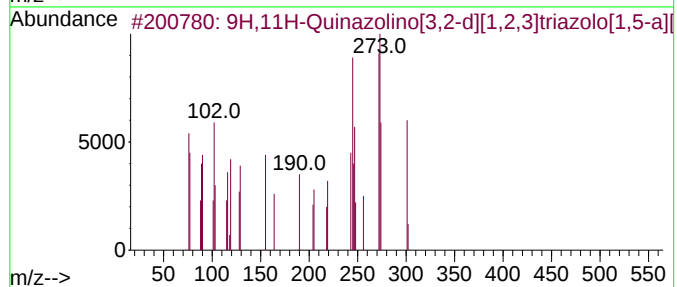
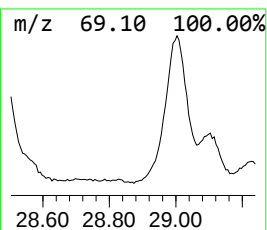
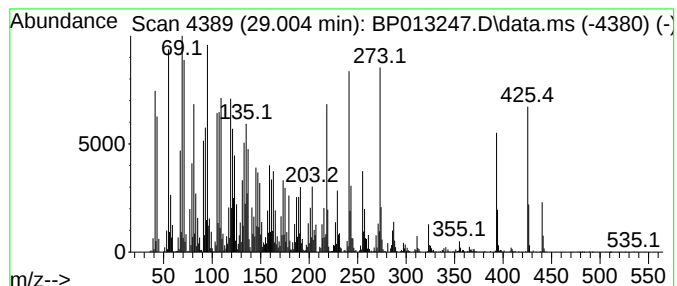
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9H,11H-Quinazolino[3,2-d][1... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
29.003	29.07 ng/ul	9409930	Perylene-d12	23.910

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9H,11H-Quinazolino[3,2-d][1,2,3]...	301	C17H11N5O	112254-02-1	91		
2	1,2,4-Oxadiazol-5(4H)-one, 4-(2-...	273	C13H8ClN3O2	288246-55-9	56		
3	D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	53		
4	Pregnan-18-ol, (5.alpha.)-	304	C21H36O	056053-09-9	49		
5	9-Thiabicyclo[3.3.1]nonane, 2,6-...	274	C14H18N4S	1000141-52-1	30		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122222\
Data File : BP013247.D
Acq On : 22 Dec 2022 15:43
Operator : CG/JU
Sample : N6015-06
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBXU3

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
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n-Hexadecanoic ...	18.216	7.7	ng/ul	1992860	4	17.345	5159220	20.0
(1H)Pyrrole, 3,...	19.828	5.6	ng/ul	1935240	5	21.433	6857380	20.0
1-Heneicosyl fo...	21.128	10.5	ng/ul	3612050	5	21.433	6857380	20.0
(DEL) Alkane: S...	23.122	4.9	ng/ul	1590790	6	23.910	6474870	20.0
(DEL) Alkane: S...	24.521	12.5	ng/ul	4048950	6	23.910	6474870	20.0
Oxirane, hexade...	25.933	8.3	ng/ul	2689290	6	23.910	6474870	20.0
(DEL) Alkane: S...	26.439	26.4	ng/ul	8529940	6	23.910	6474870	20.0
9H,11H-Quinazol...	29.003	29.1	ng/ul	9409930	6	23.910	6474870	20.0