

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
 Data File : BP013303.D
 Acq On : 24 Dec 2022 16:57
 Operator : CG/JU
 Sample : N6016-10
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBYD4

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: BP013303.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.328	21	24	32	rVB	173507	229078	1.57%	0.128%
2	3.758	94	97	112	rBV	1482926	2209987	15.16%	1.233%
3	3.864	112	115	121	rVV	135649	178525	1.22%	0.100%
4	3.940	125	128	133	rVB4	19115	26956	0.18%	0.015%
5	4.087	149	153	157	rVB	44262	54984	0.38%	0.031%
6	4.458	212	216	223	rVB5	25882	50364	0.35%	0.028%
7	4.652	246	249	257	rVB5	18234	32003	0.22%	0.018%
8	5.034	308	314	320	rVB3	34321	65510	0.45%	0.037%
9	5.117	325	328	336	rVB2	25990	43676	0.30%	0.024%
10	5.258	346	352	361	rVB	73067	131101	0.90%	0.073%
11	6.611	576	582	589	rVB	255560	394954	2.71%	0.220%
12	6.922	631	635	639	rVB4	27908	45824	0.31%	0.026%
13	7.105	658	666	679	rVB	2633968	4144999	28.43%	2.312%
14	7.269	685	694	705	rBV	2315587	3580199	24.56%	1.997%
15	7.363	708	710	721	rVB6	22389	46950	0.32%	0.026%
16	7.469	721	728	738	rBV	2944210	4557935	31.27%	2.543%
17	7.940	799	808	815	rBV	2750984	4284937	29.39%	2.390%
18	7.999	815	818	823	rVB3	16600	26009	0.18%	0.015%
19	8.652	921	929	940	rBV	2634243	4220167	28.95%	2.354%
20	9.110	997	1007	1015	rBV	2714781	4439532	30.46%	2.477%
21	9.263	1027	1033	1039	rVB2	69136	105165	0.72%	0.059%
22	9.505	1069	1074	1083	rVB	74690	126579	0.87%	0.071%
23	9.834	1122	1130	1143	rBV	2319256	3772091	25.88%	2.104%
24	9.987	1151	1156	1158	rBV4	14705	25971	0.18%	0.014%
25	10.057	1164	1168	1174	rVB6	39213	70297	0.48%	0.039%
26	10.375	1214	1222	1240	rBV	3353196	5625940	38.59%	3.138%
27	10.752	1278	1286	1300	rBV	3532231	6079165	41.70%	3.391%
28	10.893	1303	1310	1327	rVV	1774035	3169155	21.74%	1.768%
29	11.540	1412	1420	1428	rVB5	24214	49453	0.34%	0.028%
30	11.693	1436	1446	1451	rBV5	18867	56799	0.39%	0.032%
31	12.163	1522	1526	1530	rVB3	20568	32639	0.22%	0.018%
32	12.240	1535	1539	1545	rBV4	13844	25119	0.17%	0.014%
33	12.334	1550	1555	1560	rBV2	56545	93207	0.64%	0.052%
34	12.851	1637	1643	1647	rBV8	16638	28992	0.20%	0.016%
35	12.946	1655	1659	1662	rVV5	26593	47455	0.33%	0.026%
36	13.340	1718	1726	1729	rBV8	23013	57140	0.39%	0.032%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
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37	13.398	1729	1736	1739	rVV2	63073	117091	0.80%	0.065%
38	13.457	1741	1746	1751	rVV5	71500	124723	0.86%	0.070%
39	13.504	1751	1754	1758	rVV4	25731	46367	0.32%	0.026%
40	13.540	1758	1760	1769	rVV8	23451	48557	0.33%	0.027%

41	13.904	1816	1822	1828	rVB2	315691	443189	3.04%	0.247%
42	13.993	1828	1837	1845	rBV	5184647	7086377	48.61%	3.953%
43	14.051	1845	1847	1851	rVB3	25009	32537	0.22%	0.018%
44	14.104	1851	1856	1860	rVB7	40449	58884	0.40%	0.033%
45	14.234	1873	1878	1879	rBV5	20058	27943	0.19%	0.016%

46	14.281	1879	1886	1896	rBV	5643116	8455299	58.00%	4.717%
47	14.422	1903	1910	1915	rBV3	64327	100474	0.69%	0.056%
48	14.469	1915	1918	1924	rVB	52704	71280	0.49%	0.040%
49	14.540	1924	1930	1932	rBV2	83130	123823	0.85%	0.069%
50	14.587	1932	1938	1944	rVB	4764465	6830000	46.85%	3.810%

51	14.645	1944	1948	1952	rBV2	120440	168089	1.15%	0.094%
52	14.787	1966	1972	1976	rBV	1244797	1888357	12.95%	1.053%
53	14.828	1976	1979	1987	rVV2	337331	568811	3.90%	0.317%
54	14.898	1987	1991	1994	rVV	43998	81904	0.56%	0.046%
55	14.940	1994	1998	2002	rVV2	76529	124861	0.86%	0.070%

56	14.987	2002	2006	2011	rVV3	78732	142517	0.98%	0.080%
57	15.040	2011	2015	2021	rVV7	53919	93775	0.64%	0.052%
58	15.104	2024	2026	2032	rVB4	19493	28915	0.20%	0.016%
59	15.310	2058	2061	2066	rVB5	21787	30284	0.21%	0.017%
60	15.369	2066	2071	2075	rBV7	20325	41534	0.28%	0.023%

61	15.492	2084	2092	2099	rBV	218788	341931	2.35%	0.191%
62	15.581	2100	2107	2113	rBV	6821979	9650119	66.20%	5.383%
63	15.628	2113	2115	2121	rVB	32983	43092	0.30%	0.024%
64	15.698	2121	2127	2136	rBV	1299684	1735215	11.90%	0.968%
65	15.822	2143	2148	2154	rVB4	62294	105963	0.73%	0.059%

66	15.887	2154	2159	2160	rBV4	21754	27459	0.19%	0.015%
67	15.945	2164	2169	2178	rVB	534297	745066	5.11%	0.416%
68	16.028	2178	2183	2184	rBV	103855	137518	0.94%	0.077%
69	16.045	2184	2186	2193	rVB4	124320	192374	1.32%	0.107%
70	16.151	2198	2204	2210	rBV2	229568	326720	2.24%	0.182%

71	16.210	2210	2214	2221	rBV8	43145	71836	0.49%	0.040%
72	16.292	2224	2228	2234	rVB8	31926	69664	0.48%	0.039%
73	16.475	2256	2259	2262	rVV4	25278	37306	0.26%	0.021%
74	16.510	2262	2265	2270	rVB2	57254	78040	0.54%	0.044%
75	16.728	2298	2302	2308	rBV2	80828	124772	0.86%	0.070%

76	17.222	2380	2386	2389	rBV3	37872	67261	0.46%	0.038%
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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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 Peak separation: 5

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77	17.345	2400	2407	2417	rVV	4379134	6389097	43.83%	3.564%
78	17.445	2417	2424	2431	rVB	5279578	7774971	53.34%	4.337%
79	17.698	2463	2467	2470	rBV3	54313	83018	0.57%	0.046%
80	17.998	2516	2518	2522	rVB4	29298	31224	0.21%	0.017%
81	18.104	2531	2536	2539	rBV5	48620	88107	0.60%	0.049%
82	18.157	2541	2545	2550	rBV	145065	218733	1.50%	0.122%
83	18.216	2550	2555	2565	rBV	1825579	2565058	17.60%	1.431%
84	18.545	2608	2611	2621	rVB2	113563	178307	1.22%	0.099%
85	19.204	2720	2723	2727	rBV2	62188	77993	0.54%	0.044%
86	19.251	2729	2731	2736	rVB3	74644	82405	0.57%	0.046%
87	19.333	2737	2745	2753	rBV3	898120	2035448	13.96%	1.135%
88	19.445	2760	2764	2776	rVB	547601	777307	5.33%	0.434%
89	19.675	2798	2803	2815	rVB	7280459	9783978	67.12%	5.458%
90	19.822	2825	2828	2835	rVB	195267	231935	1.59%	0.129%
91	20.975	3022	3024	3028	rBV	167835	165812	1.14%	0.092%
92	21.127	3046	3050	3057	rBV	2000721	2882602	19.77%	1.608%
93	21.433	3097	3102	3107	rBV	5527023	7344928	50.39%	4.097%
94	22.069	3207	3210	3220	rVB	1100282	1746679	11.98%	0.974%
95	23.127	3385	3390	3396	rBV	2031192	3273365	22.46%	1.826%
96	23.757	3490	3497	3510	rVB2	5482698	12029369	82.52%	6.711%
97	23.916	3518	3524	3540	rVB2	4005674	8144561	55.87%	4.543%
98	24.533	3622	3629	3637	rVB	3242556	6691797	45.91%	3.733%
99	24.633	3641	3646	3656	rVB	1376132	3467053	23.78%	1.934%
100	28.498	4295	4303	4323	rVB2	3792681	14577167	100.00%	8.132%

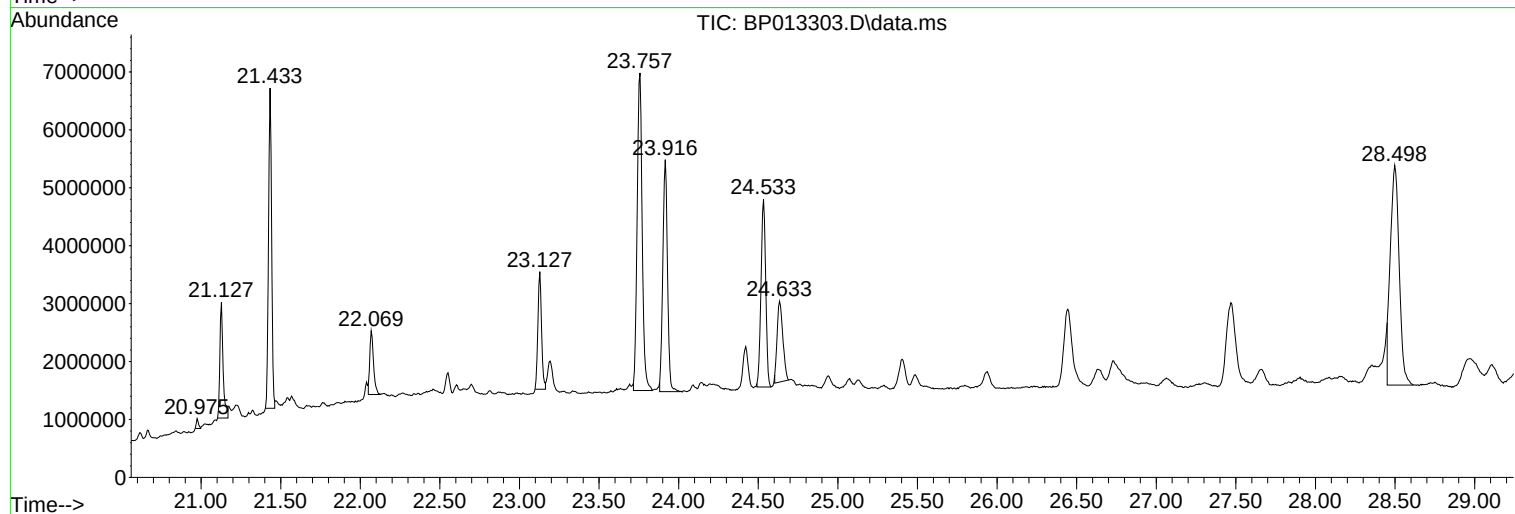
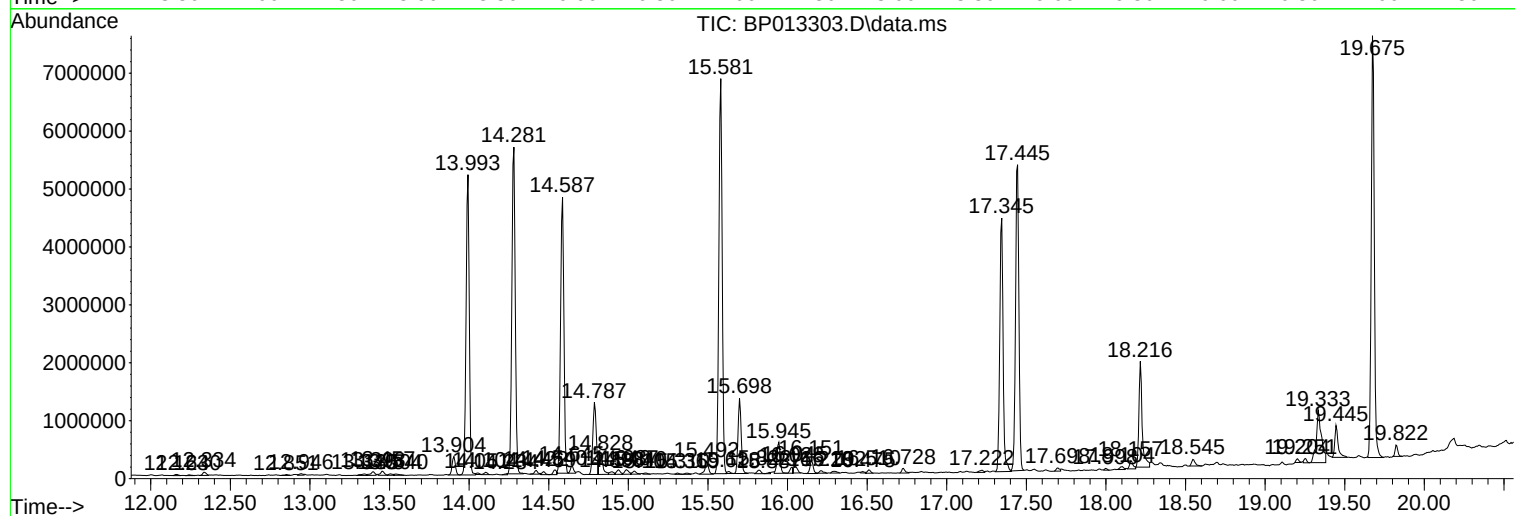
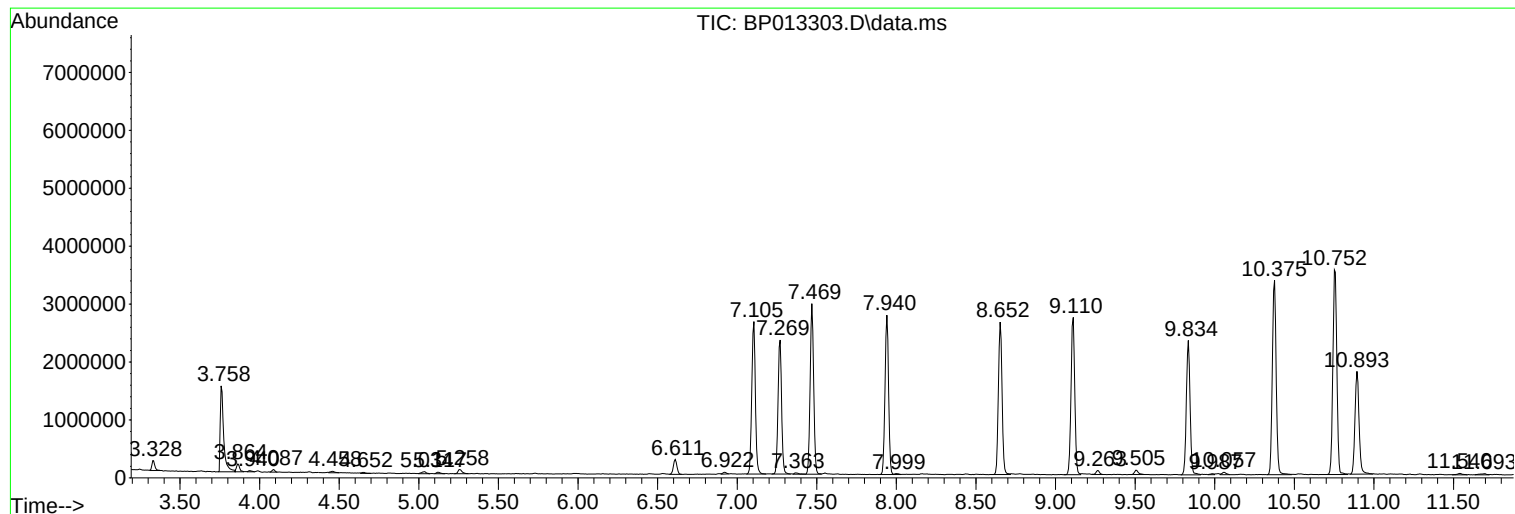
Sum of corrected areas: 179259698

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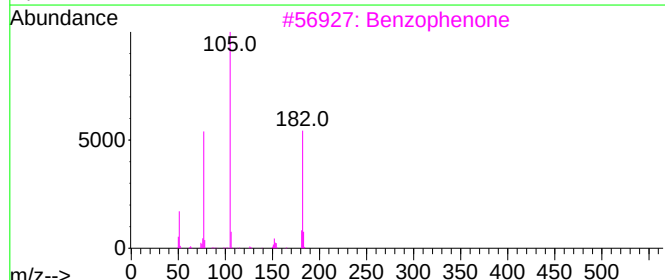
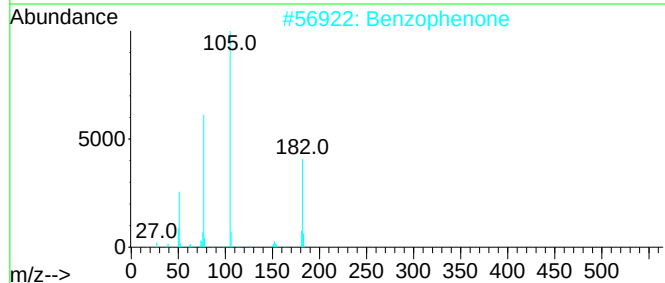
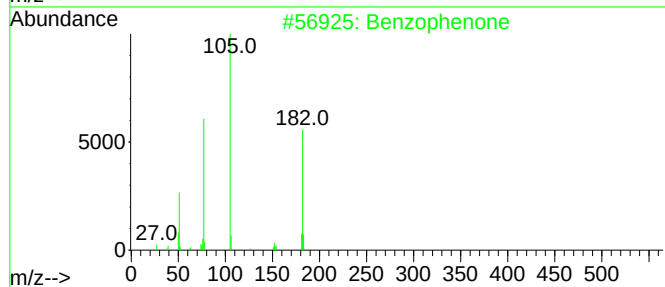
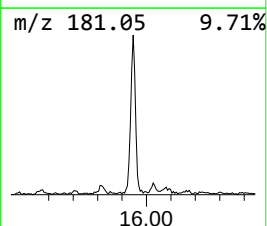
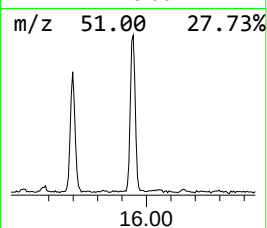
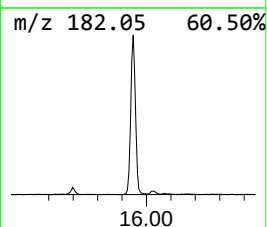
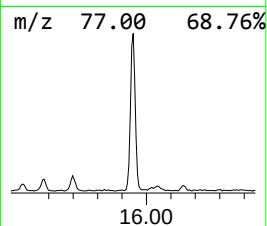
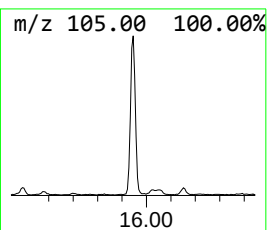
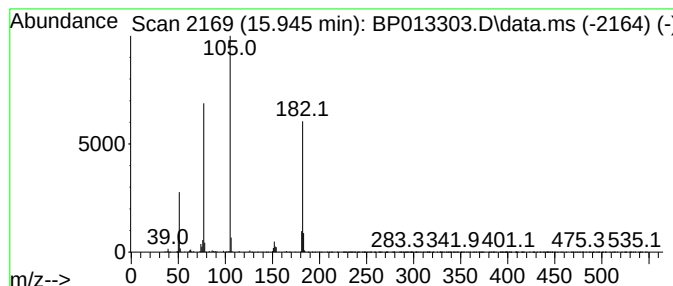
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzophenone Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.945	2.18 ng/ul	745066	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	96
3			Benzophenone	182	C13H10O	000119-61-9	95
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzenamine, N-(3-pyridinylmethy...	182	C12H10N2	029722-97-2	42



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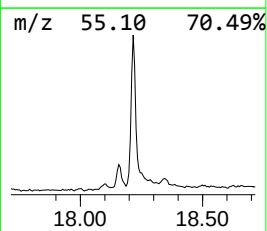
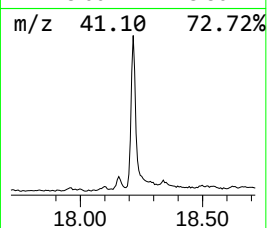
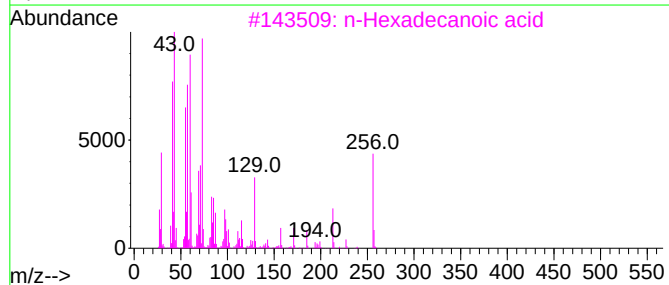
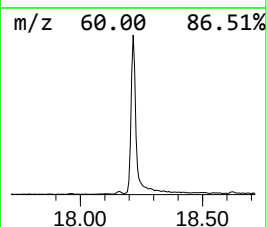
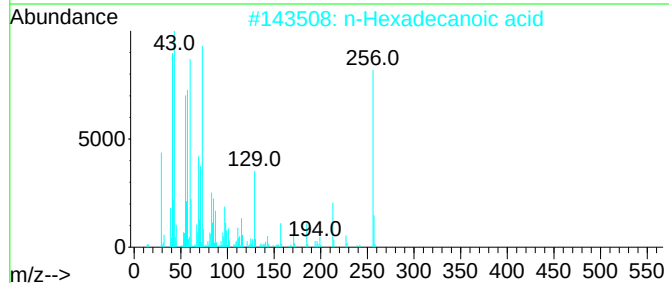
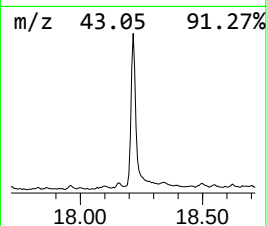
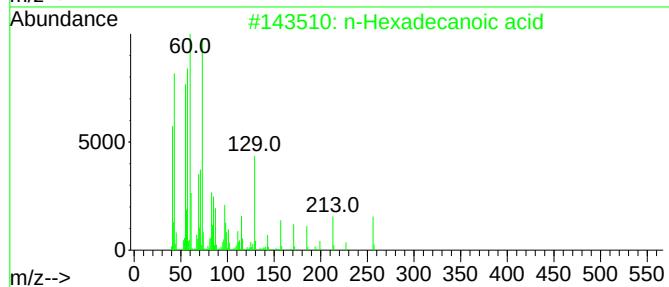
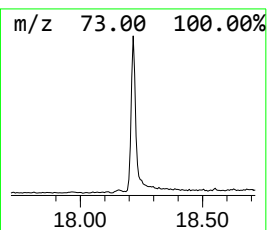
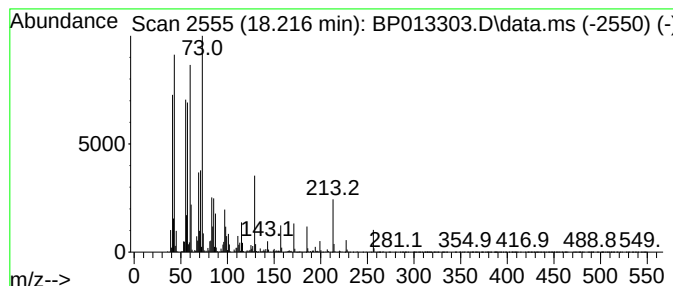
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 n-Hexadecanoic acid Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.216	8.03 ng/ul	2565060	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	97
5			Tridecanoic acid	214	C13H26O2	000638-53-9	95



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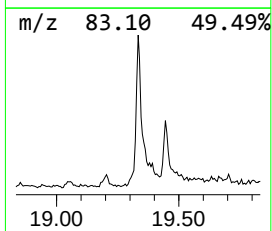
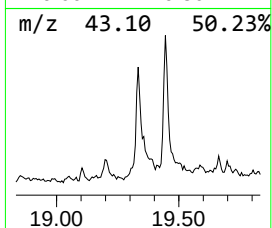
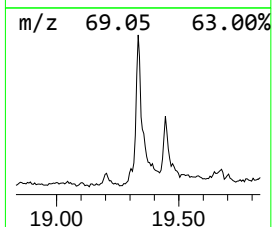
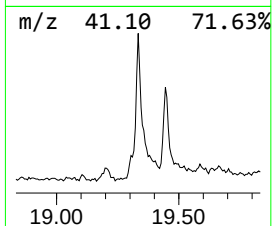
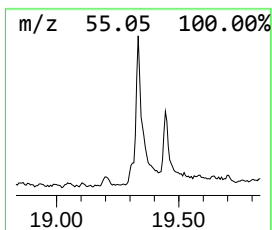
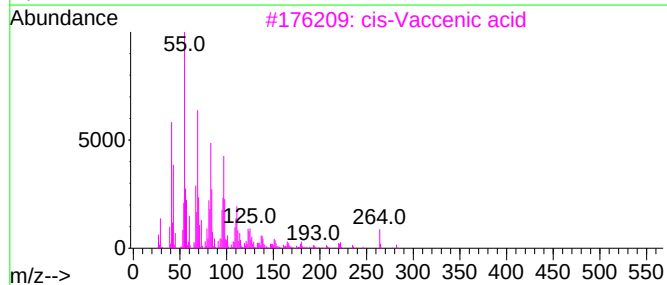
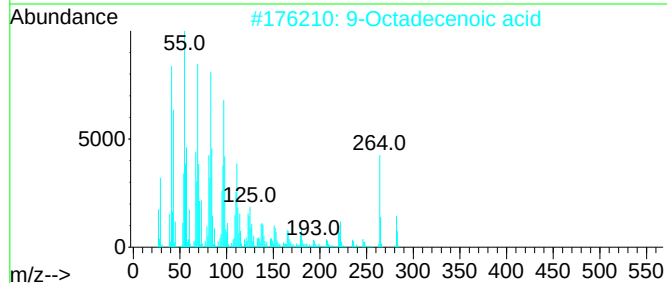
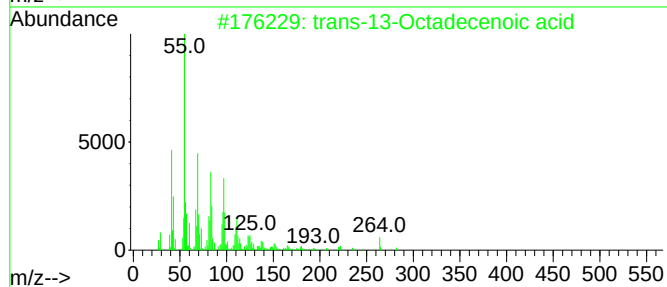
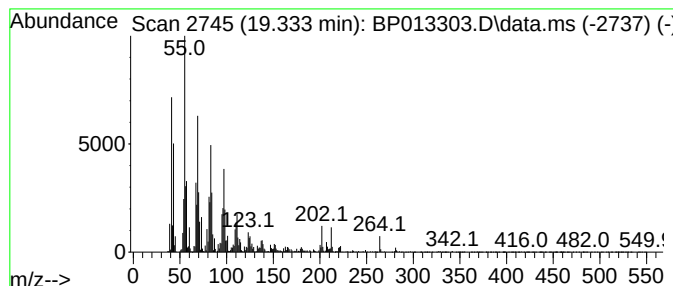
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP120722.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 trans-13-Octadecenoic acid Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.333	6.37 ng/ul	2035450	Phenanthrene-d10	17.345

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			trans-13-Octadecenoic acid	282	C18H34O2	000693-71-0	97
2			9-Octadecenoic acid	282	C18H34O2	002027-47-6	97
3			cis-Vaccenic acid	282	C18H34O2	000506-17-2	97
4			cis-13-Octadecenoic acid	282	C18H34O2	013126-39-1	96
5			Oleyl alcohol, trifluoroacetate	364	C20H35F3O2	1000352-68-4	93



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Data File : BP013303.D
Acq On : 24 Dec 2022 16:57
Operator : CG/JU
Sample : N6016-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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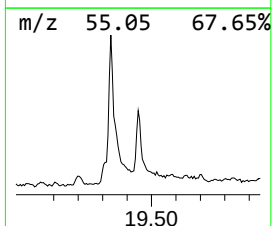
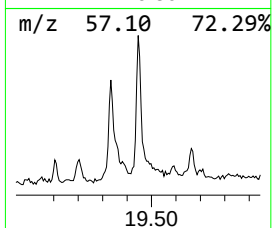
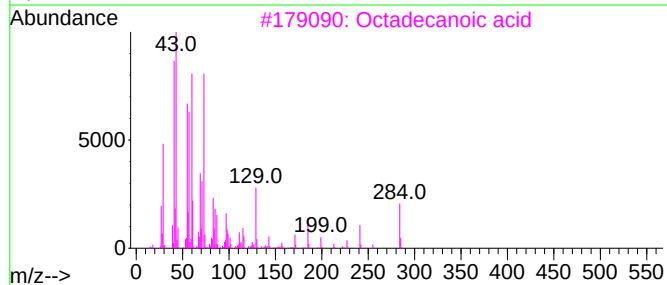
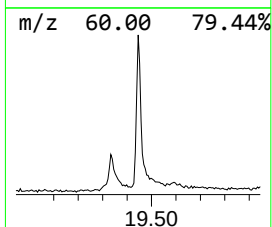
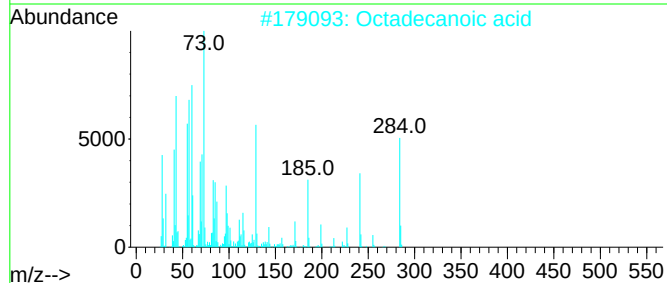
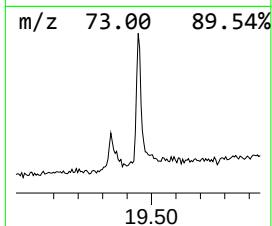
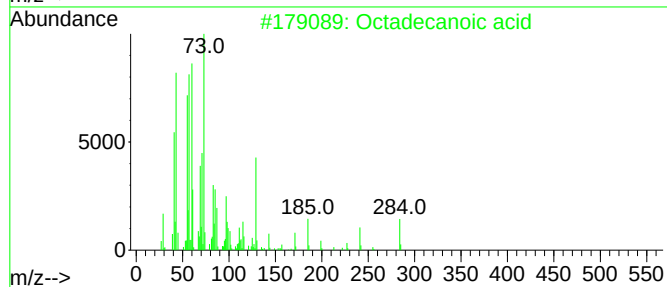
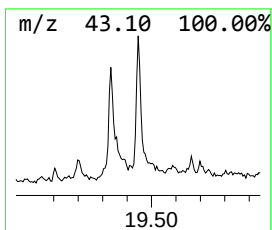
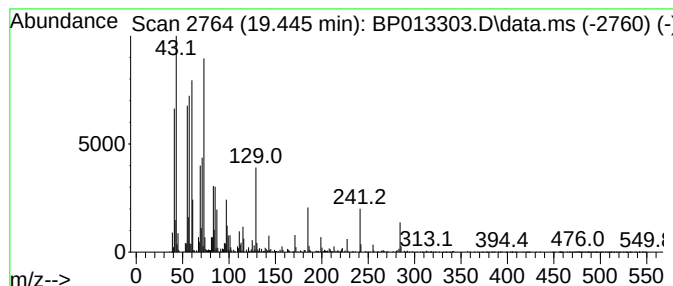
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	2.12 ng/ul	777307	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octadecanoic acid	284	C18H36O2	000057-11-4	99
2			Octadecanoic acid	284	C18H36O2	000057-11-4	99
3			Octadecanoic acid	284	C18H36O2	000057-11-4	98
4			Octadecanoic acid	284	C18H36O2	000057-11-4	94
5			Octadecanoic acid	284	C18H36O2	000057-11-4	93



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Misc :
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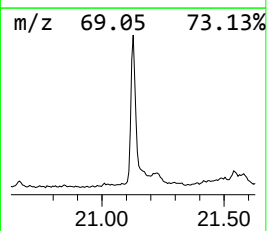
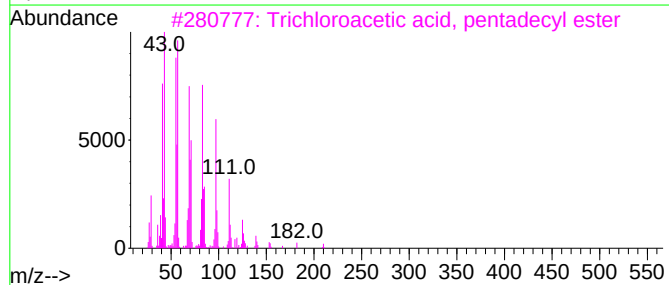
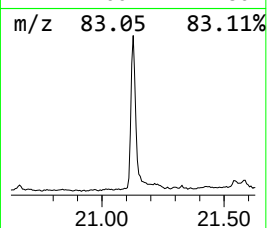
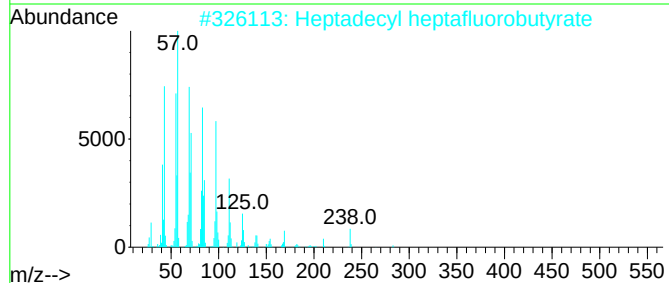
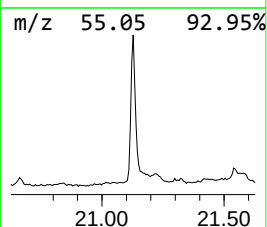
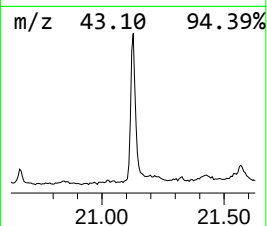
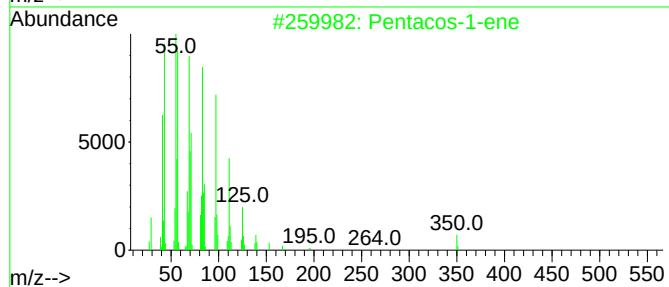
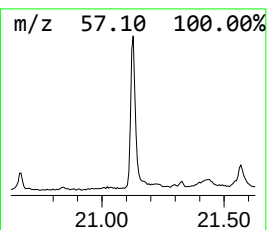
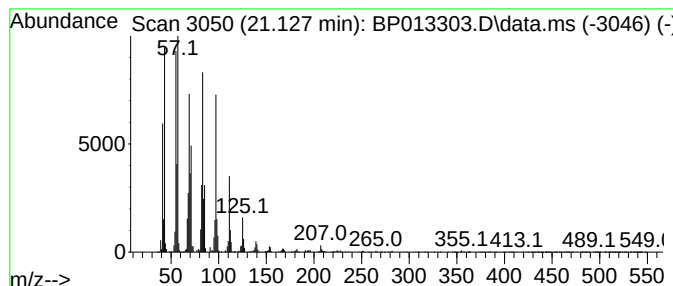
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.127	7.85 ng/ul	2882600	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Pentacos-1-ene	350	C25H50	016980-85-1	94
2			Heptadecyl heptafluorobutyrate	452	C21H35F7O2	959085-66-6	94
3			Trichloroacetic acid, pentadecyl...	372	C17H31Cl3O2	074339-53-0	94
4			Pentafluoropropionic acid, penta...	374	C18H31F5O2	959092-08-1	94
5			Trichloroacetic acid, hexadecyl ...	386	C18H33Cl3O2	074339-54-1	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013303.D
Acq On : 24 Dec 2022 16:57
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Sample : N6016-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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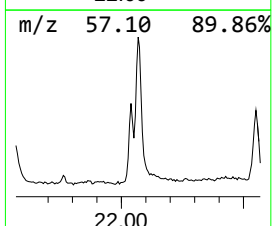
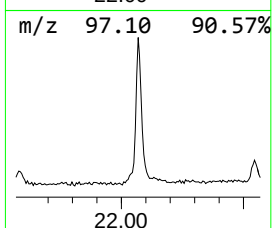
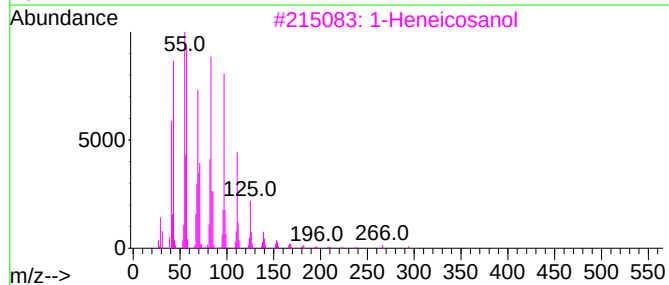
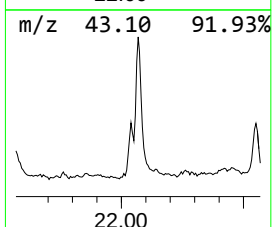
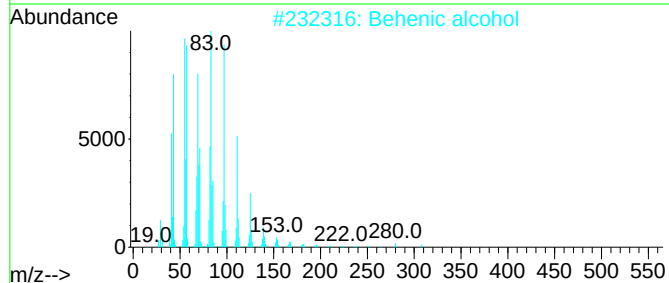
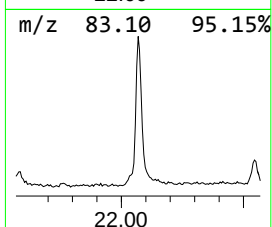
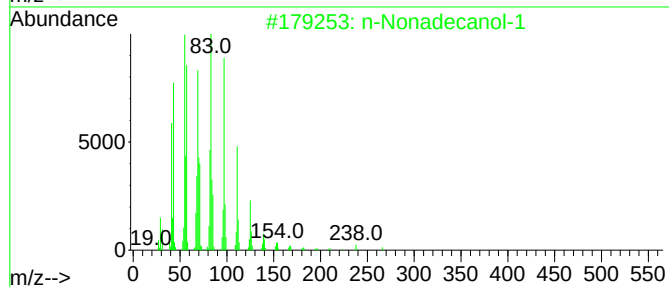
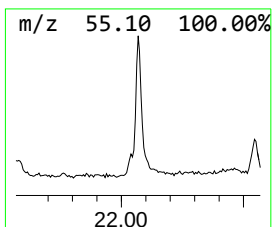
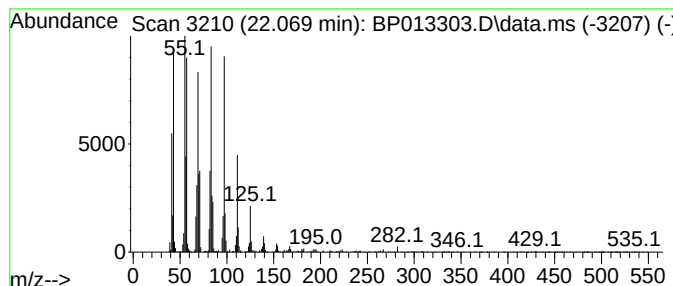
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 n-Nonadecanol-1 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.069	4.76 ng/ul	1746680	Chrysene-d12	21.433

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			n-Nonadecanol-1	284	C19H40O	001454-84-8	94
2			Behenic alcohol	326	C22H46O	000661-19-8	94
3			1-Heneicosanol	312	C21H44O	015594-90-8	94
4			1-Nonadecene	266	C19H38	018435-45-5	93
5			1-Hexacosene	364	C26H52	018835-33-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013303.D
Acq On : 24 Dec 2022 16:57
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Sample : N6016-10
Misc :
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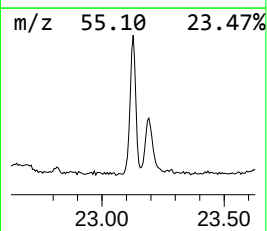
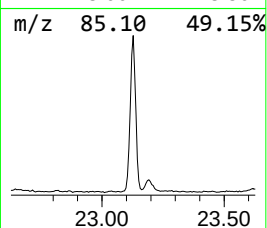
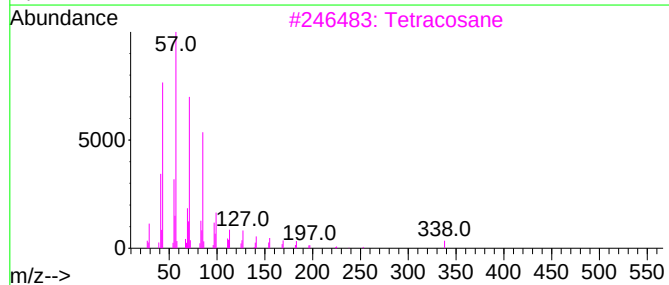
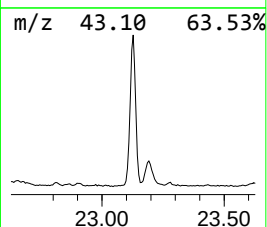
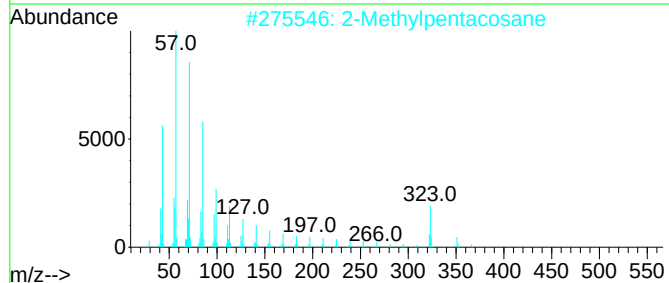
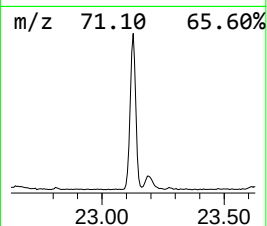
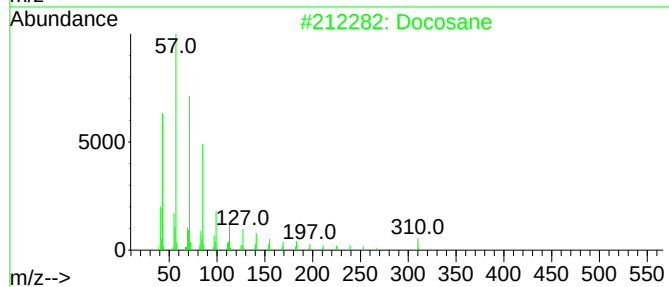
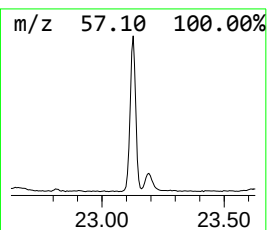
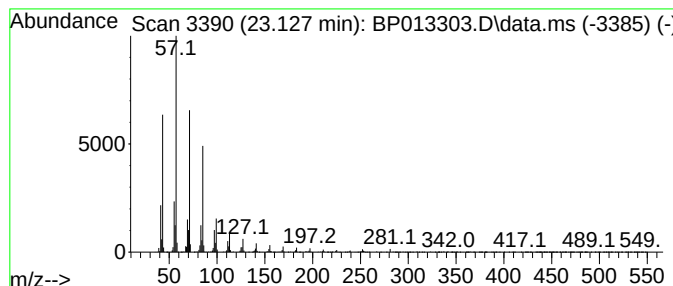
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 (DEL) Alkane: Straight-Chai... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
23.127	8.04 ng/ul	3273370	Perylene-d12	23.916

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Docosane	310	C22H46	000629-97-0	95
2		2-Methylpentacosane	366	C26H54	000629-87-8	94
3		Tetracosane	338	C24H50	000646-31-1	91
4		Octacosane	394	C28H58	000630-02-4	91
5		Heneicosane	296	C21H44	000629-94-7	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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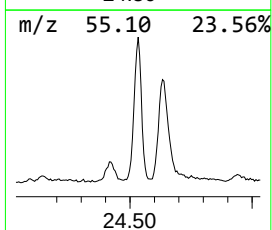
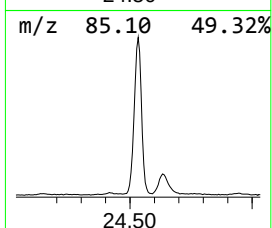
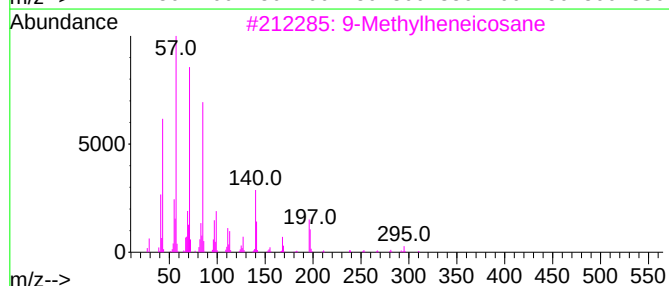
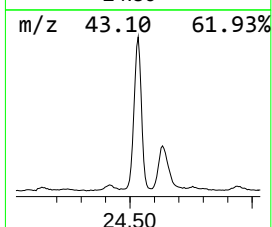
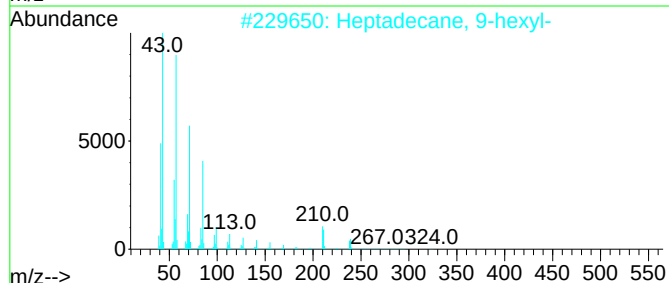
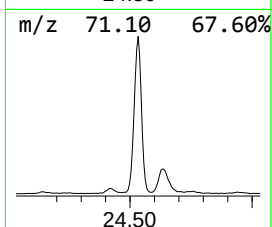
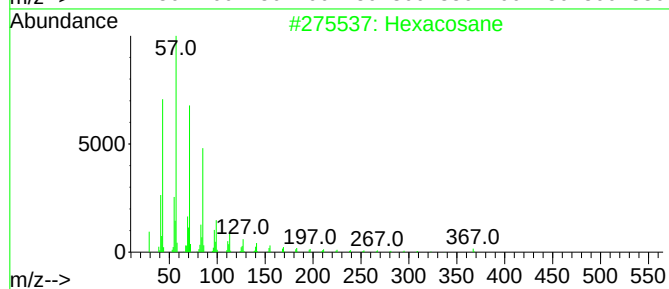
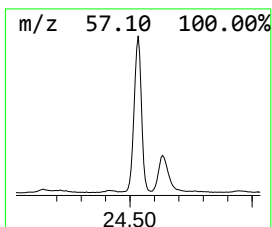
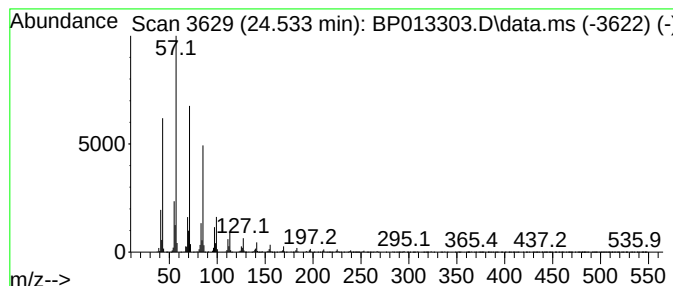
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD		R.T.	
24.533	16.43 ng/ul	6691800	Perylene-d12		23.916	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Hexacosane		366	C26H54	000630-01-3	94
2	Heptadecane, 9-hexyl-		324	C23H48	055124-79-3	94
3	9-Methylheneicosane		310	C22H46	070475-51-3	91
4	Tetratetracontane		619	C44H90	007098-22-8	91
5	Heptadecane, 9-octyl-		352	C25H52	007225-64-1	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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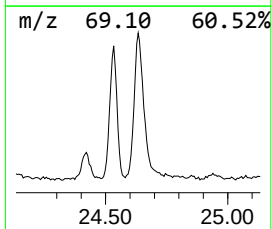
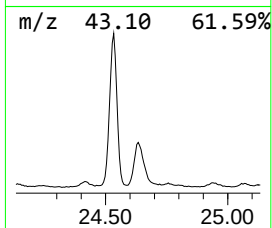
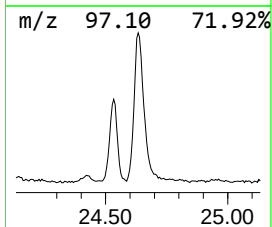
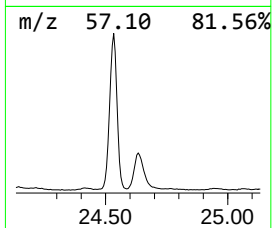
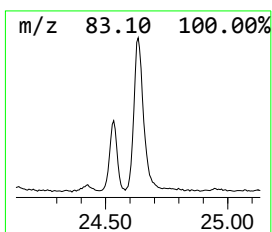
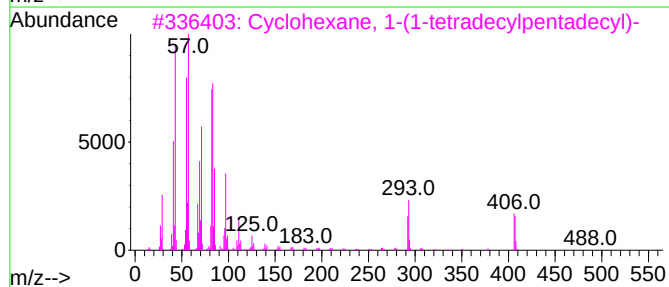
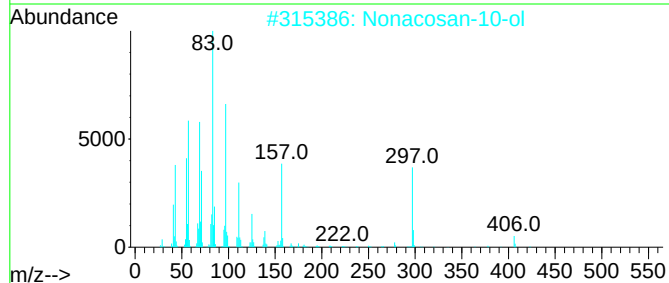
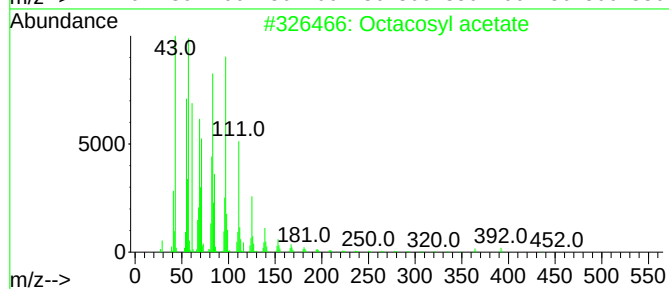
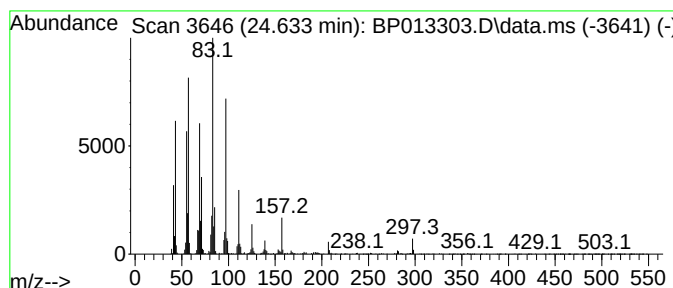
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Octacosyl acetate Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
24.633	8.51 ng/ul	3467050	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Octacosyl acetate	452	C30H60O2	018206-97-8	95
2			Nonacosan-10-ol	424	C29H60O	000504-55-2	91
3			Cyclohexane, 1-(1-tetradecylpent...	491	C35H70	055521-27-2	49
4			Octacosyl trifluoroacetate	506	C30H57F3O2	1000351-74-9	49
5			Tetracosyl pentafluoropropionate	500	C27H49F5O2	1000351-81-3	49



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
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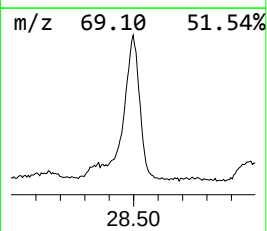
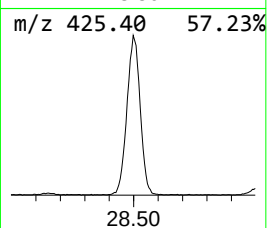
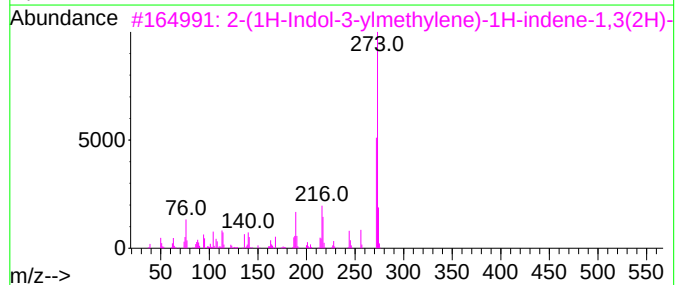
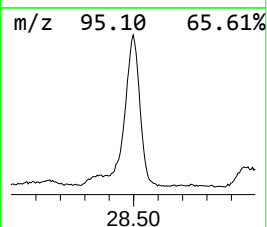
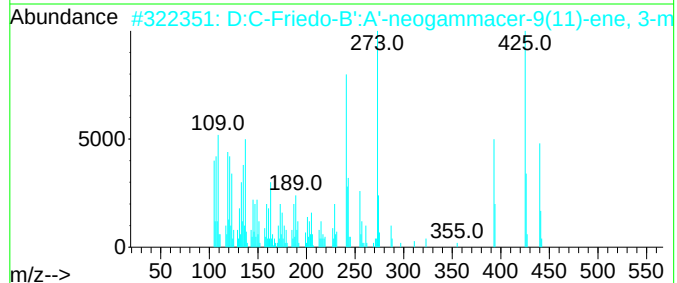
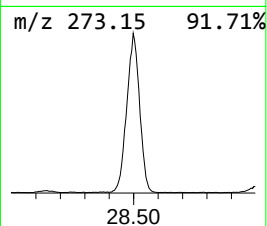
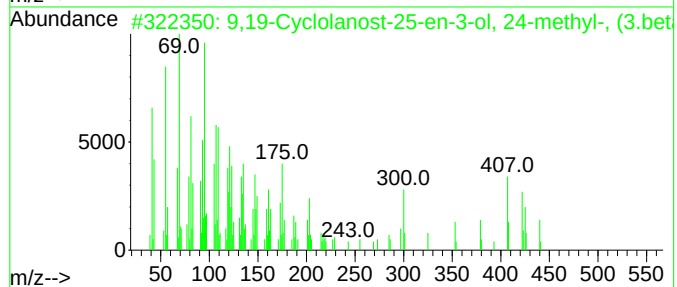
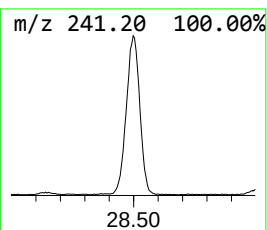
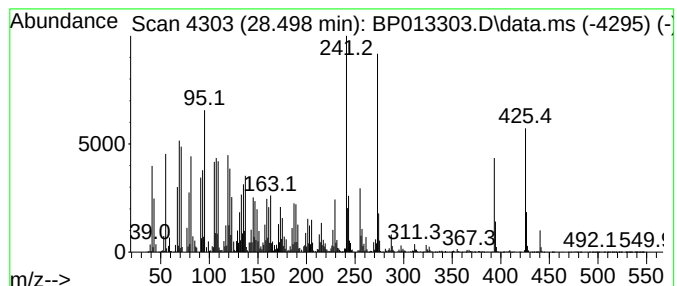
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 unknown-01 Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
28.498	35.80 ng/ul	14577200	Perylene-d12	23.916

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	9,19-Cyclolanost-25-en-3-ol, 24-...	440	C31H52O	000511-61-5	46		
2	D:C-Friedo-B':A'-neogammacer-9(1...	440	C31H52O	004555-56-0	46		
3	2-(1H-Indol-3-ylmethylene)-1H-in...	273	C18H11NO2	031060-64-7	25		
4	Lupeol	426	C30H50O	000545-47-1	25		
5	Imidazolo[1,2-a]pyridine, 6-chlo...	273	C13H8ClN3O2	118000-62-7	25		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP122422\
Data File : BP013303.D
Acq On : 24 Dec 2022 16:57
Operator : CG/JU
Sample : N6016-10
Misc :
ALS Vial : 14 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBYD4

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	2.2	ng/ul	745066	3	14.587	6830000	20.0
n-Hexadecanoic ...	18.216	8.0	ng/ul	2565060	4	17.345	6389100	20.0
trans-13-Octade...	19.333	6.4	ng/ul	2035450	4	17.345	6389100	20.0
Octadecanoic acid	19.445	2.1	ng/ul	777307	5	21.433	7344930	20.0
(DEL) Alkane: S...	21.127	7.8	ng/ul	2882600	5	21.433	7344930	20.0
n-Nonadecanol-1	22.069	4.8	ng/ul	1746680	5	21.433	7344930	20.0
(DEL) Alkane: S...	23.127	8.0	ng/ul	3273370	6	23.916	8144560	20.0
(DEL) Alkane: S...	24.533	16.4	ng/ul	6691800	6	23.916	8144560	20.0
Octacosyl acetate	24.633	8.5	ng/ul	3467050	6	23.916	8144560	20.0
unknown-01	28.498	35.8	ng/ul	14577200	6	23.916	8144560	20.0