

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
 Data File : BP013754.D
 Acq On : 03 Feb 2023 22:55
 Operator : CG/JU
 Sample : 01381-14
 Misc :
 ALS Vial : 55 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 A44N3

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-BP020123.M
 Title : SVOA CALIBRATION

Signal : TIC: BP013754.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.440	38	43	57	rBV2	304502	605730	5.59%	0.593%
2	3.917	121	124	142	rBV	132336	263074	2.43%	0.258%
3	4.146	158	163	169	rVB2	17160	31590	0.29%	0.031%
4	4.246	176	180	185	rBV	14127	21842	0.20%	0.021%
5	5.199	338	342	354	rVB	38516	64128	0.59%	0.063%
6	5.405	369	377	390	rBV	1836624	3039479	28.04%	2.976%
7	5.599	406	410	417	rBV2	16284	29658	0.27%	0.029%
8	6.775	604	610	618	rBV	42140	69962	0.65%	0.068%
9	7.099	658	665	671	rBV5	8624	18173	0.17%	0.018%
10	7.275	689	695	706	rVB	150947	263358	2.43%	0.258%
11	7.446	718	724	733	rBV	791415	1262793	11.65%	1.236%
12	7.658	753	760	769	rBV	650790	1085449	10.01%	1.063%
13	8.016	812	821	829	rBV	130200	217842	2.01%	0.213%
14	8.128	834	840	843	rVV	689129	1128738	10.41%	1.105%
15	8.163	843	846	857	rVB	933300	1496997	13.81%	1.466%
16	8.487	893	901	914	rBV	4508482	7449250	68.73%	7.293%
17	8.834	953	960	969	rBV	505644	840178	7.75%	0.823%
18	8.940	972	978	986	rVV	51042	95506	0.88%	0.094%
19	9.028	988	993	1001	rVV6	14901	34724	0.32%	0.034%
20	9.122	1005	1009	1014	rVB2	14237	22030	0.20%	0.022%
21	9.299	1031	1039	1052	rBV	804773	1401049	12.93%	1.372%
22	9.657	1096	1100	1104	rBV6	6189	11469	0.11%	0.011%
23	9.705	1104	1108	1113	rVV	23733	35362	0.33%	0.035%
24	9.763	1113	1118	1126	rVB	28918	49503	0.46%	0.048%
25	9.975	1147	1154	1157	rBV2	41315	72898	0.67%	0.071%
26	10.028	1157	1163	1178	rVB	673979	1110330	10.24%	1.087%
27	10.569	1246	1255	1265	rBV	1176789	1964409	18.12%	1.923%
28	10.816	1290	1297	1312	rVB	908203	1506421	13.90%	1.475%
29	10.957	1313	1321	1333	rBV	1064741	1770373	16.33%	1.733%
30	11.093	1334	1344	1359	rBV	784989	1444659	13.33%	1.414%
31	11.346	1379	1387	1396	rVB	156266	279398	2.58%	0.274%
32	11.622	1429	1434	1442	rBV2	7002	13953	0.13%	0.014%
33	12.269	1537	1544	1549	rBV8	5557	11127	0.10%	0.011%
34	12.463	1569	1577	1584	rBV	176336	289982	2.68%	0.284%
35	12.528	1584	1588	1600	rVB2	26012	58564	0.54%	0.057%
36	12.975	1658	1664	1677	rVB	112680	188122	1.74%	0.184%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
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37	13.287	1711	1717	1726	rBV	5988	14526	0.13%	0.014%
38	13.581	1758	1767	1771	rBV5	16122	33362	0.31%	0.033%
39	13.651	1774	1779	1785	rVB2	30520	53287	0.49%	0.052%
40	13.822	1803	1808	1815	rVB3	13952	23051	0.21%	0.023%
41	14.169	1859	1867	1882	rBV	2522138	3720247	34.32%	3.642%
42	14.287	1882	1887	1892	rVB8	13892	21855	0.20%	0.021%
43	14.345	1892	1897	1900	rVB2	9487	12103	0.11%	0.012%
44	14.463	1910	1917	1933	rVB	2852510	4365533	40.28%	4.274%
45	14.640	1944	1947	1951	rVB5	8595	10907	0.10%	0.011%
46	14.763	1960	1968	1986	rBV	1780501	2791144	25.75%	2.733%
47	14.951	1994	2000	2012	rBV2	133720	259542	2.39%	0.254%
48	15.157	2032	2035	2042	rVB8	11021	20842	0.19%	0.020%
49	15.257	2047	2052	2057	rBV4	13686	25254	0.23%	0.025%
50	15.757	2128	2137	2149	rBV	4239036	6130966	56.56%	6.003%
51	15.863	2149	2155	2169	rVV	963128	1331109	12.28%	1.303%
52	15.992	2172	2177	2187	rVB3	44985	113142	1.04%	0.111%
53	16.116	2192	2198	2208	rBV	1056371	1448942	13.37%	1.419%
54	16.198	2208	2212	2216	rVV2	30449	39543	0.36%	0.039%
55	16.251	2216	2221	2229	rVV	252308	376671	3.48%	0.369%
56	16.345	2231	2237	2246	rVV	48030	82306	0.76%	0.081%
57	16.534	2267	2269	2274	rVB6	8079	11396	0.11%	0.011%
58	16.681	2290	2294	2303	rBV	46232	70530	0.65%	0.069%
59	16.892	2326	2330	2336	rVB9	13462	21259	0.20%	0.021%
60	17.163	2372	2376	2386	rVB3	32977	76486	0.71%	0.075%
61	17.263	2391	2393	2397	rVV5	9097	12466	0.12%	0.012%
62	17.304	2398	2400	2405	rVB4	9279	11190	0.10%	0.011%
63	17.386	2410	2414	2423	rBV	30966	53650	0.49%	0.053%
64	17.516	2430	2436	2446	rBV	2689186	3884033	35.83%	3.803%
65	17.616	2447	2453	2463	rBV	4991946	7196625	66.40%	7.046%
66	17.763	2474	2478	2480	rBV5	12231	17640	0.16%	0.017%
67	17.798	2480	2484	2493	rVB4	42535	79352	0.73%	0.078%
68	17.881	2495	2498	2503	rVV6	15411	22342	0.21%	0.022%
69	18.251	2558	2561	2566	rVB6	14048	17383	0.16%	0.017%
70	18.369	2576	2581	2590	rBV	317501	466535	4.30%	0.457%
71	18.692	2632	2636	2646	rVB	111940	160219	1.48%	0.157%
72	19.198	2717	2722	2732	rBV2	267835	478938	4.42%	0.469%
73	19.410	2753	2758	2761	rBV2	320178	454335	4.19%	0.445%
74	19.445	2761	2764	2783	rVB	1685000	2445443	22.56%	2.394%
75	19.592	2786	2789	2796	rBV4	131296	225851	2.08%	0.221%
76	19.833	2824	2830	2837	rBV	7428142	10011730	92.37%	9.802%

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77	19.922	2837	2845	2858	rVB2	1719194	3206836	29.59%	3.140%
78	20.092	2870	2874	2881	rBV2	156096	231285	2.13%	0.226%
79	21.010	3027	3030	3035	rBV	183294	215221	1.99%	0.211%
80	21.269	3070	3074	3083	rBV2	808311	1245512	11.49%	1.219%
81	21.598	3124	3130	3136	rBV	3657730	5102555	47.08%	4.996%
82	22.174	3225	3228	3232	rVB2	196267	232827	2.15%	0.228%
83	22.898	3347	3351	3359	rVB	609215	876344	8.09%	0.858%
84	23.998	3530	3538	3555	rVB2	5071380	10838799	100.00%	10.612%
85	24.163	3559	3566	3576	rVB2	2482880	5349718	49.36%	5.238%

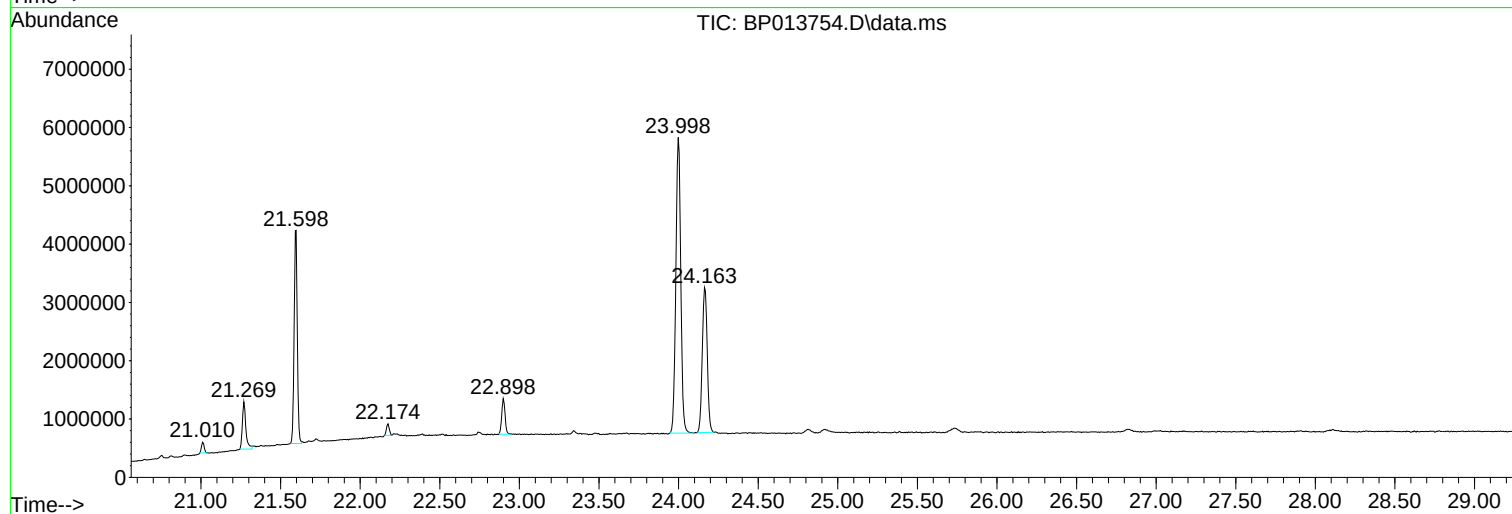
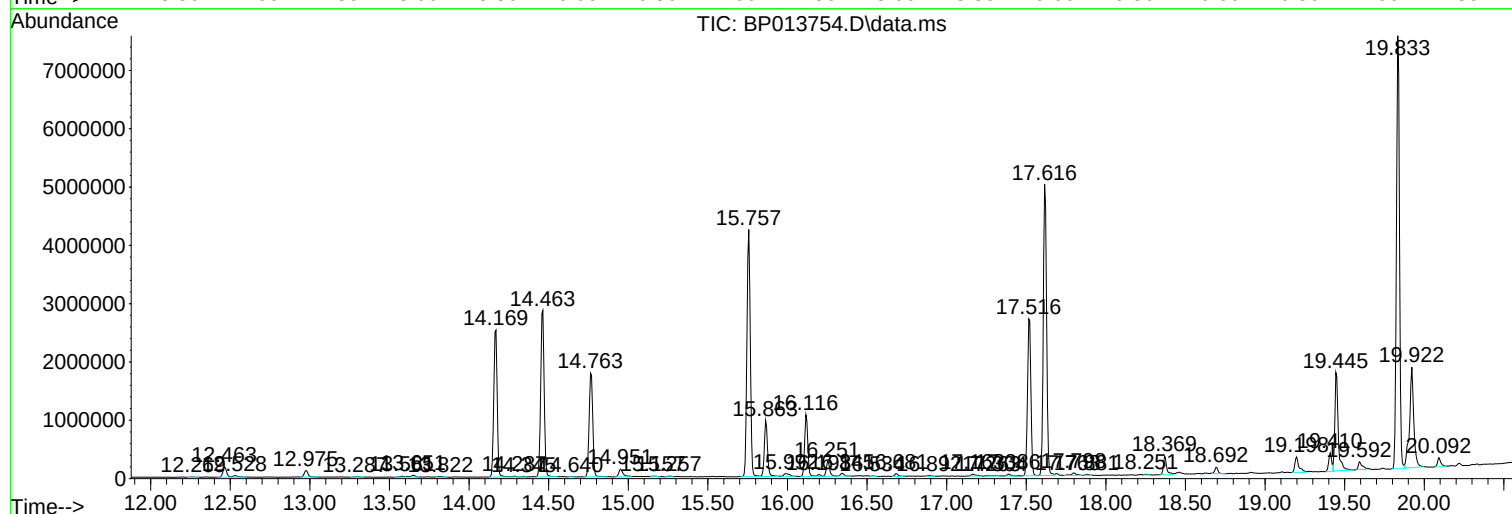
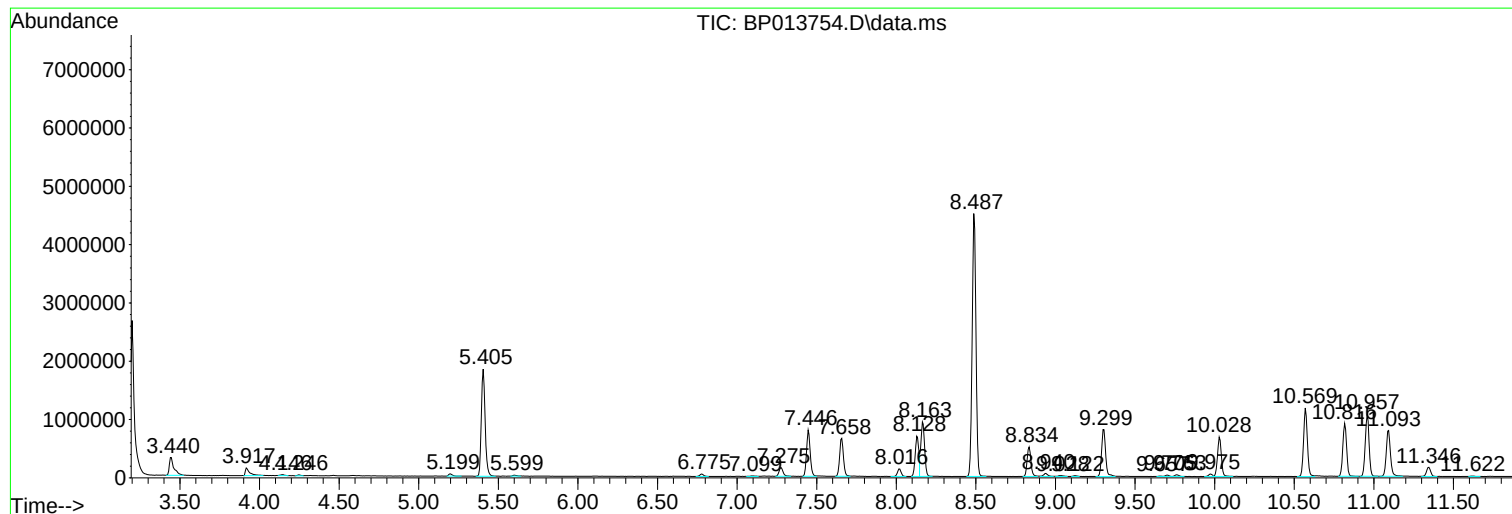
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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



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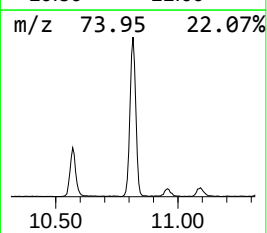
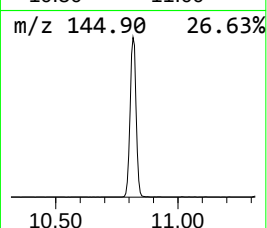
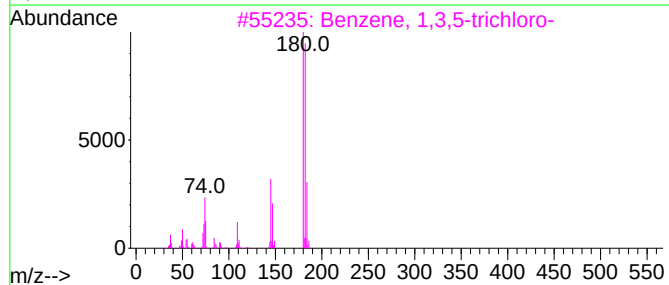
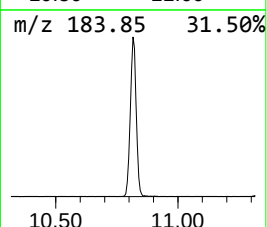
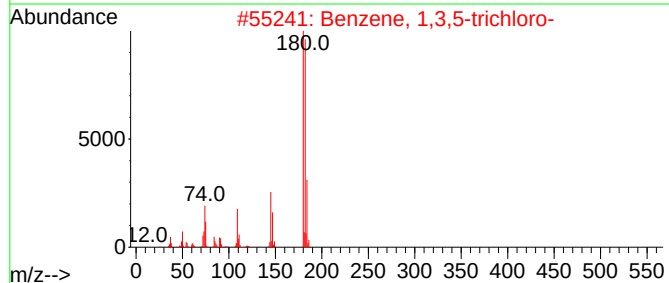
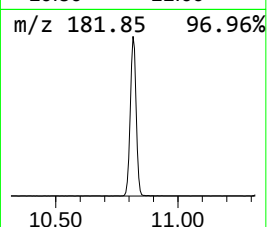
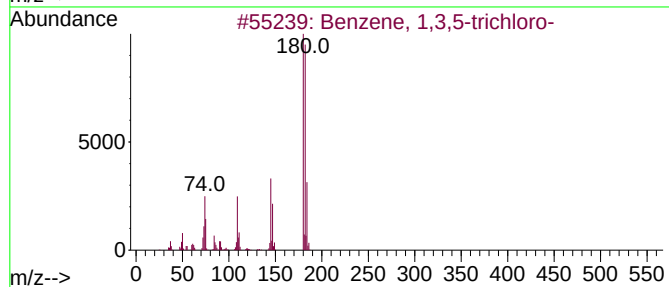
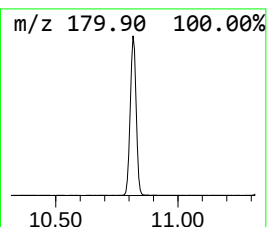
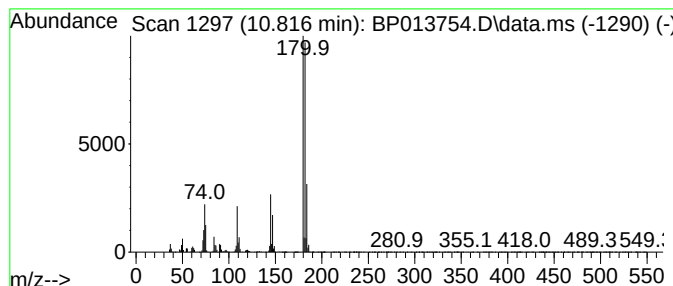
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,3,5-trichloro- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.816	17.02 ng/ul	1506420	Naphthalene-d8	10.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
2			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
3			Benzene, 1,3,5-trichloro-	180	C6H3Cl3	000108-70-3	98
4			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	97
5			Benzene, 1,2,4-trichloro-	180	C6H3Cl3	000120-82-1	95



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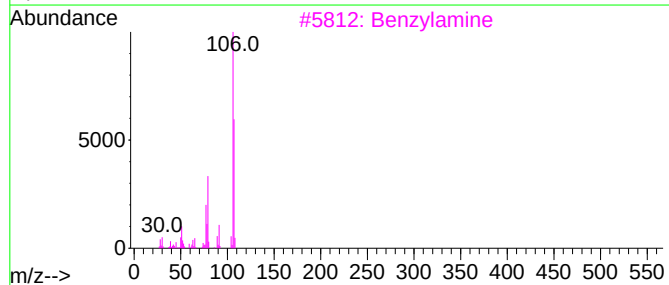
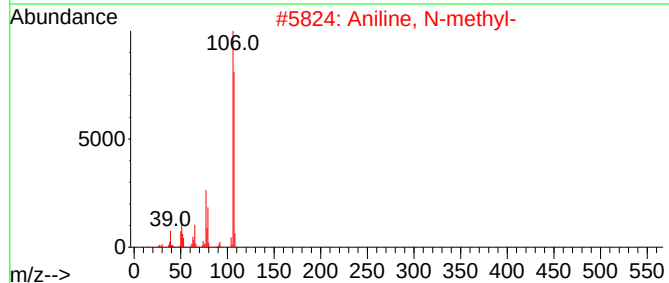
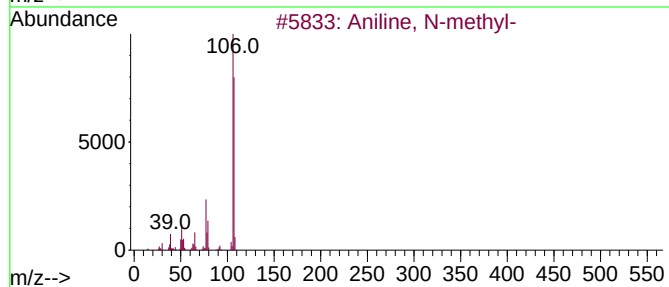
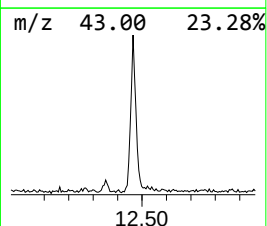
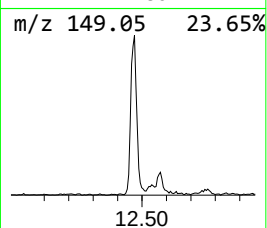
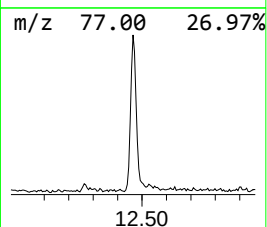
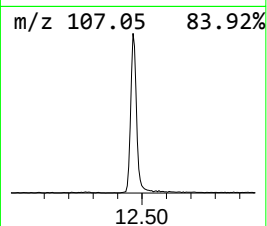
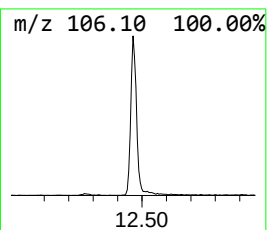
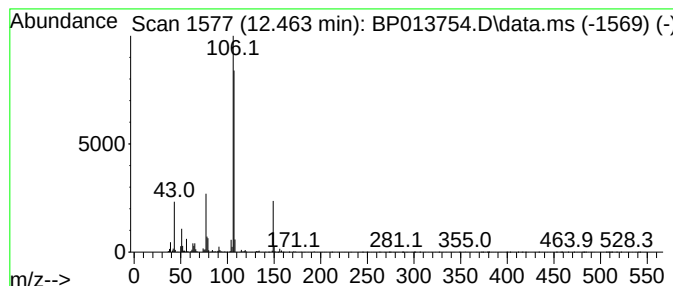
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Aniline, N-methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.463	3.28 ng/ul	289982	Naphthalene-d8	10.957

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Aniline, N-methyl-		107	C7H9N	000100-61-8	87	
2	Aniline, N-methyl-		107	C7H9N	000100-61-8	83	
3	Benzylamine		107	C7H9N	000100-46-9	80	
4	Aniline, N-methyl-		107	C7H9N	000100-61-8	80	
5	Acetamide, N-methyl-N-phenyl-		149	C9H11NO	000579-10-2	76	



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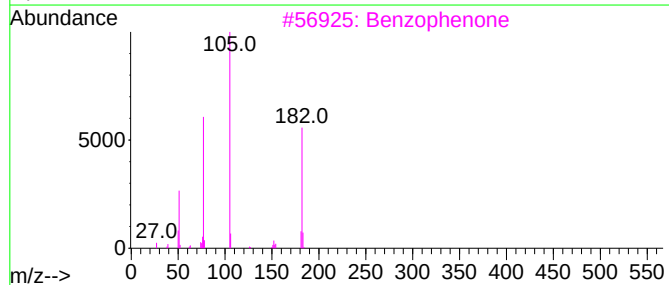
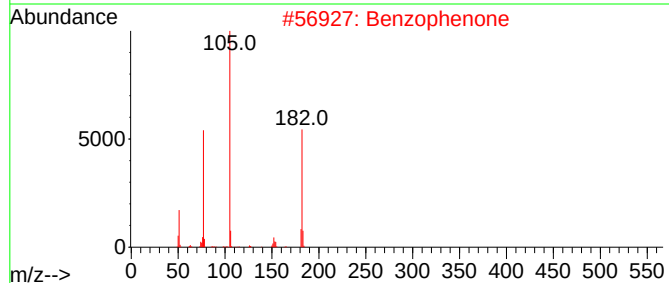
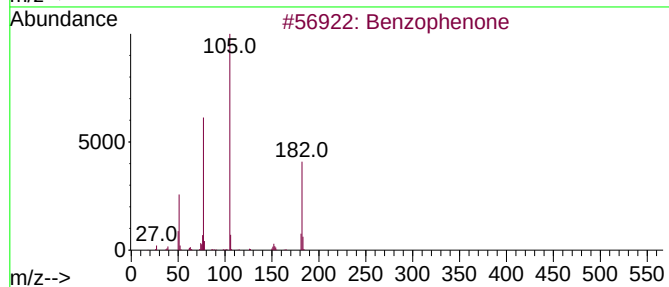
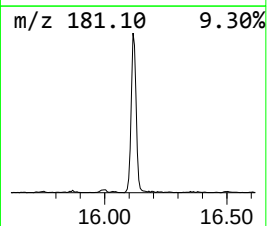
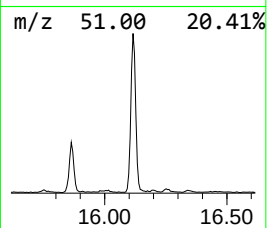
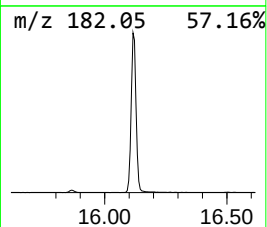
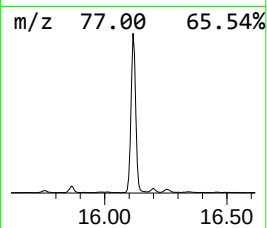
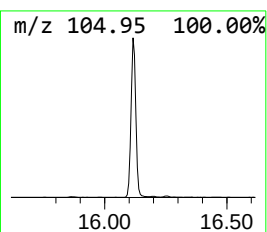
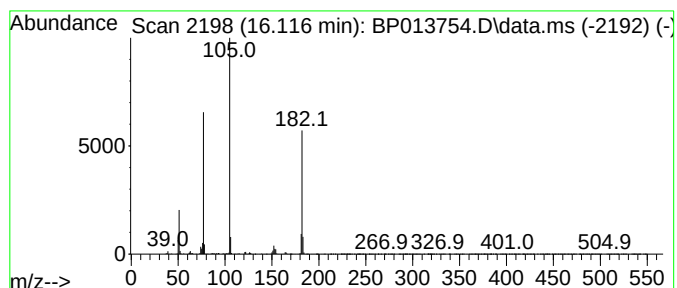
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzophenone Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.116	10.38 ng/ul	1448940	Acenaphthene-d10	14.763

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	96
4			Benzophenone	182	C13H10O	000119-61-9	95
5			1,2,4,5-Tetroxane, 3,3,6,6-tetra...	396	C26H20O4	016204-36-7	64



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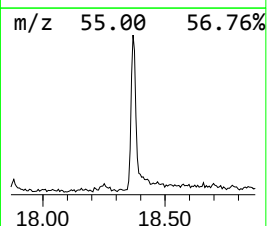
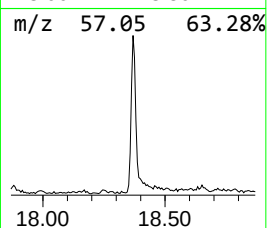
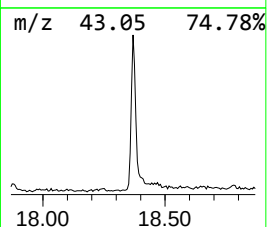
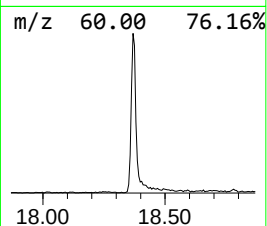
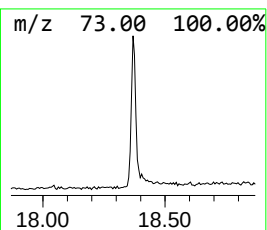
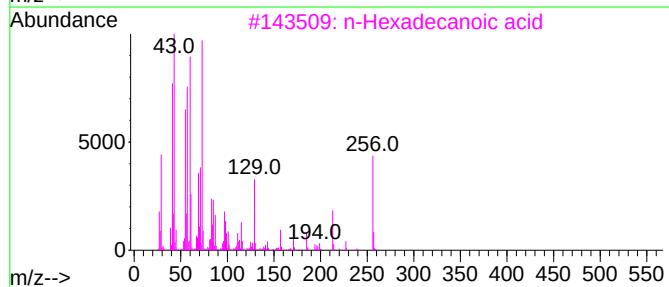
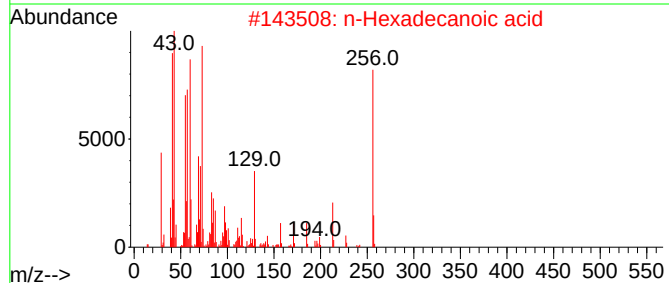
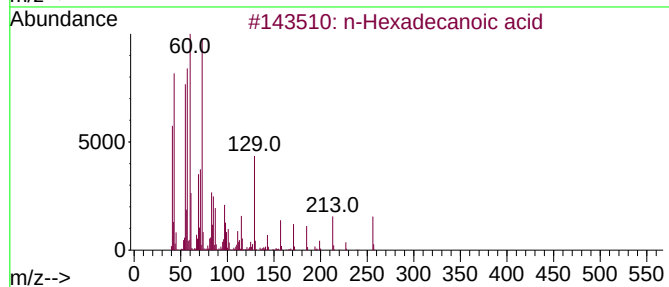
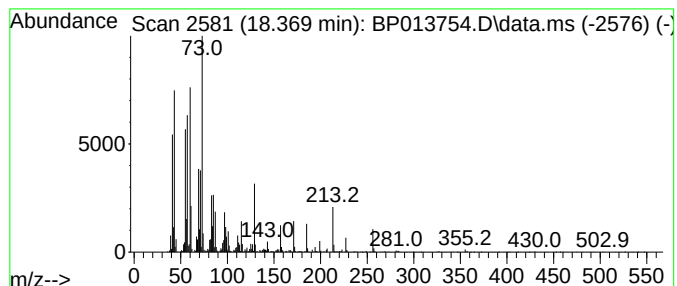
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ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.	
18.369	2.40 ng/ul	466535	Phenanthrene-d10	17.516	
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	91
5	Tridecanoic acid	214	C13H26O2	000638-53-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
Operator : CG/JU
Sample : 01381-14
Misc :
ALS Vial : 55 Sample Multiplier: 1

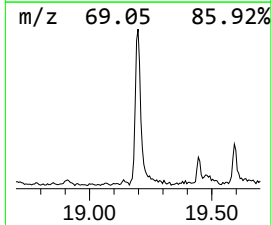
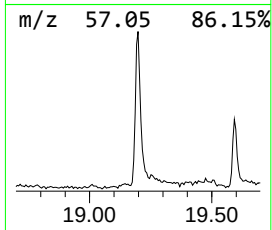
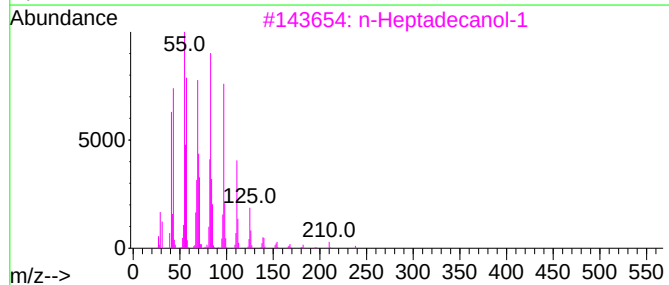
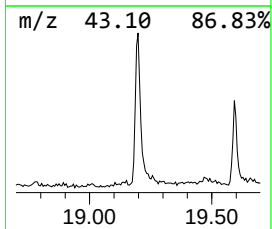
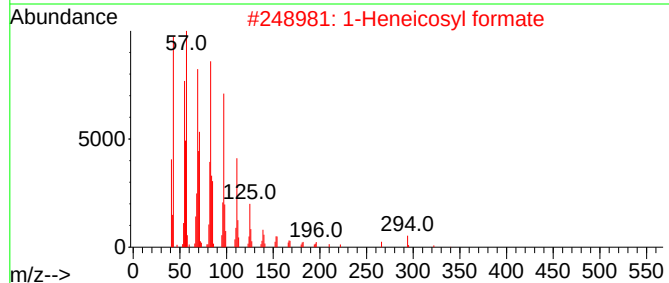
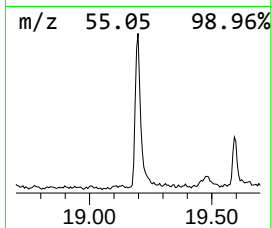
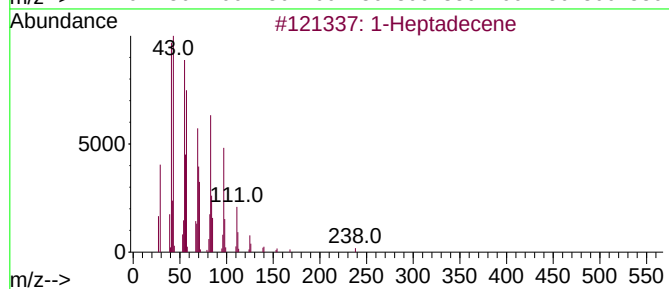
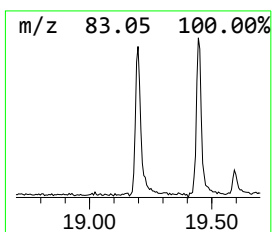
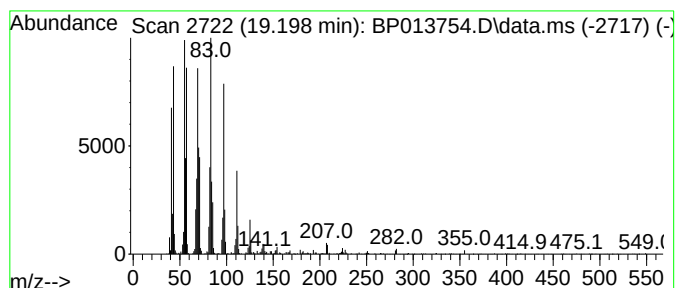
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.M
Quant Title : SVOA CALIBRATION

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

Peak Number 9 (DEL) Alkane: Straight-Chai... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD		R.T.	
19.198	2.47 ng/ul	478938	Phenanthrene-d10		17.516	
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	1-Heptadecene		238	C17H34	006765-39-5	99
2	1-Heneicosyl formate		340	C22H44O2	077899-03-7	96
3	n-Heptadecanol-1		256	C17H36O	001454-85-9	94
4	n-Nonadecanol-1		284	C19H40O	001454-84-8	94
5	Bromoacetic acid, hexadecyl ester		362	C18H35BrO2	005454-48-8	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
Operator : CG/JU
Sample : 01381-14
Misc :
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Instrument :
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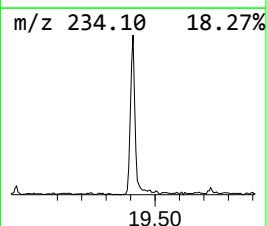
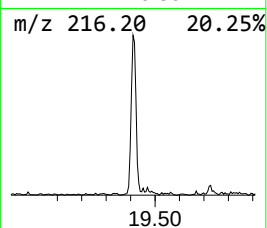
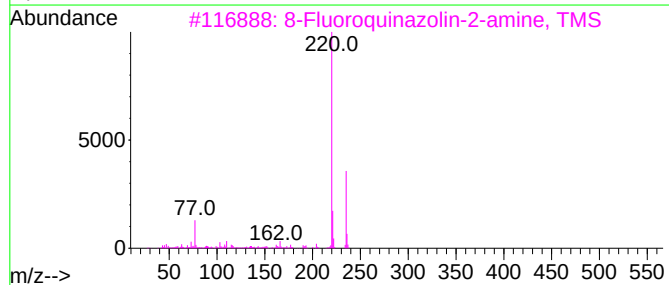
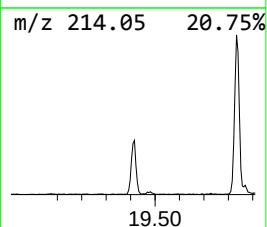
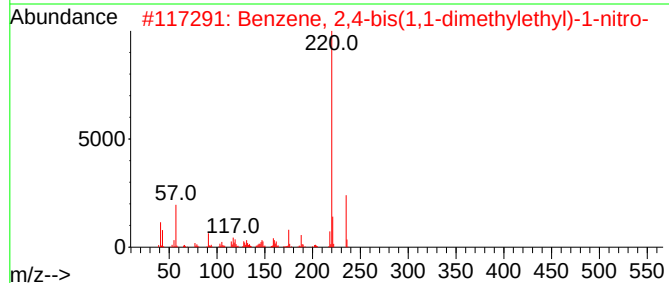
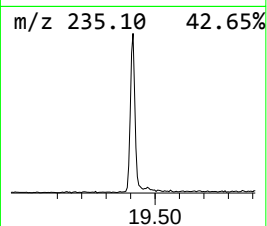
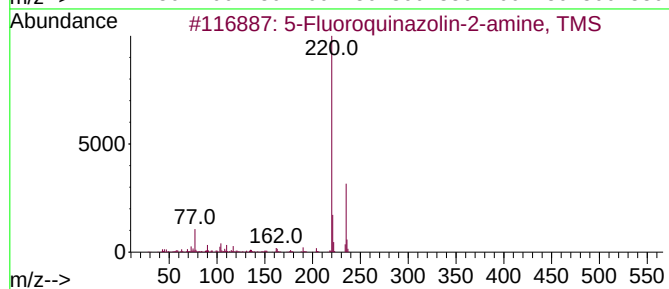
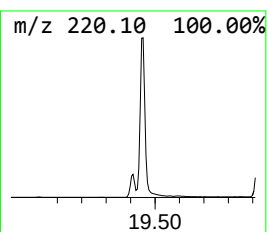
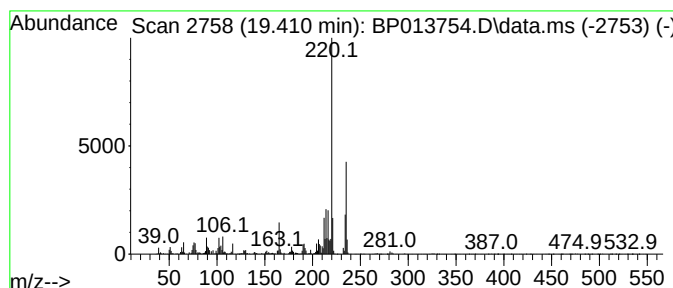
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 5-Fluoroquinazolin-2-amine,... Concentration Rank 14

R.T.	EstConc	Area	Relative to ISTD		R.T.
19.410	2.34 ng/ul	454335	Phenanthrene-d10		17.516

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	5-Fluoroquinazolin-2-amine, TMS	235	C11H14FN3Si	1000497-20-4	58	
2	Benzene, 2,4-bis(1,1-dimethyleth...	235	C14H21NO2	093720-37-7	50	
3	8-Fluoroquinazolin-2-amine, TMS	235	C11H14FN3Si	1000497-20-9	50	
4	1,5-Diphenylpyrazole	220	C15H12N2	006831-89-6	42	
5	Phenanthrene, 2,3,5-trimethyl-	220	C17H16	003674-73-5	38	



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
Operator : CG/JU
Sample : 01381-14
Misc :
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Instrument :
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ClientSampleId :
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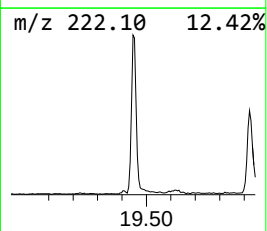
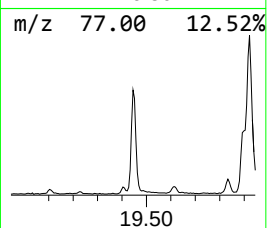
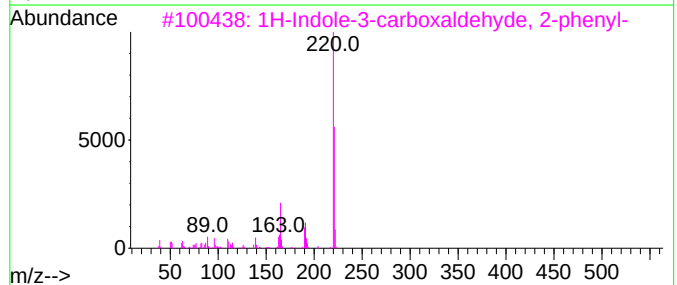
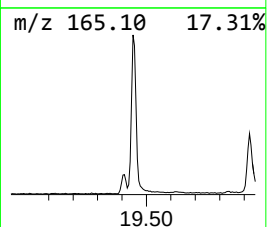
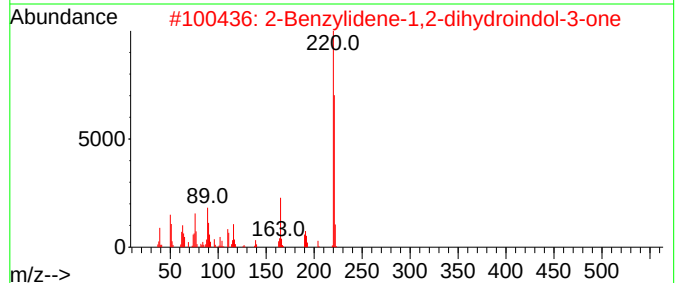
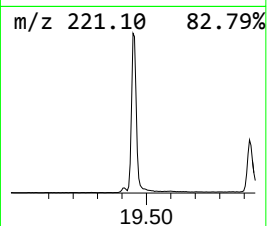
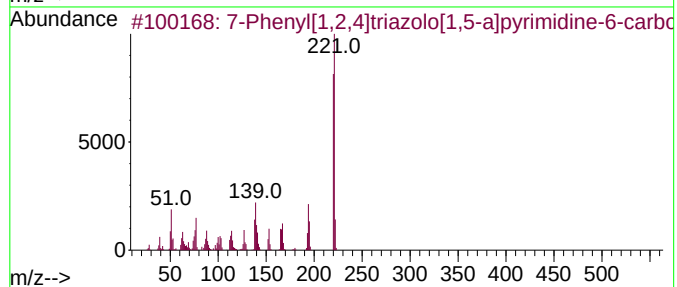
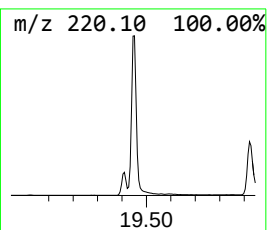
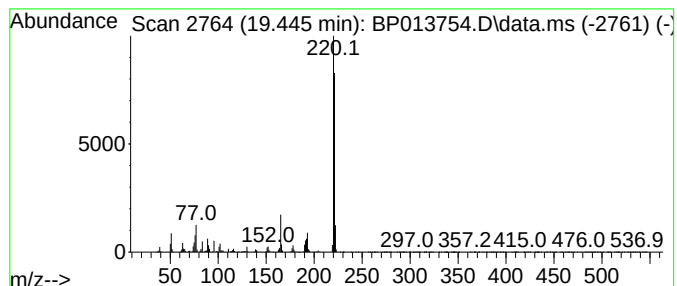
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 7-Phenyl[1,2,4]triazolo[1,5... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.445	12.59 ng/ul	2445440	Phenanthrene-d10	17.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Phenyl[1,2,4]triazolo[1,5-a]py...	221	C12H7N5	264927-73-3	86
2			2-Benzylidene-1,2-dihydroindol-3...	221	C15H11NO	039830-75-6	83
3			1H-Indole-3-carboxaldehyde, 2-ph...	221	C15H11NO	025365-71-3	78
4			1H-Quinolin-2-one, 3-phenyl-	221	C15H11NO	038035-81-3	68
5			N,N-Dimethyl-4-tert-butyl-thiobe...	221	C13H19NS	072864-57-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
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Misc :
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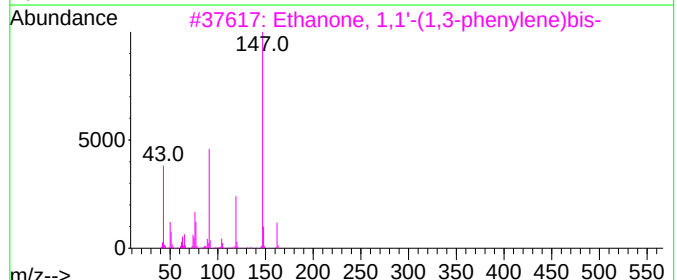
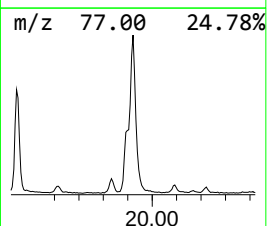
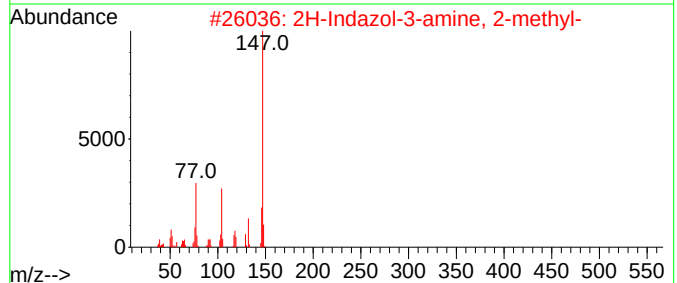
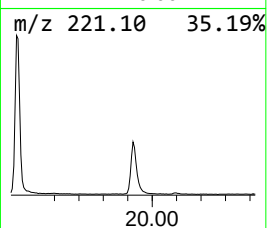
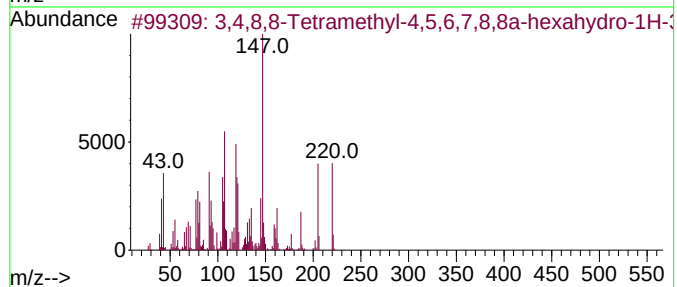
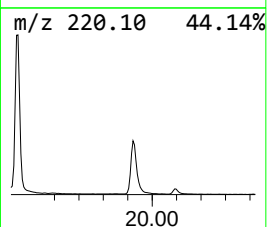
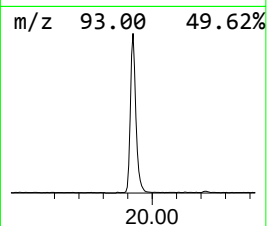
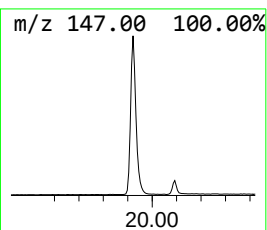
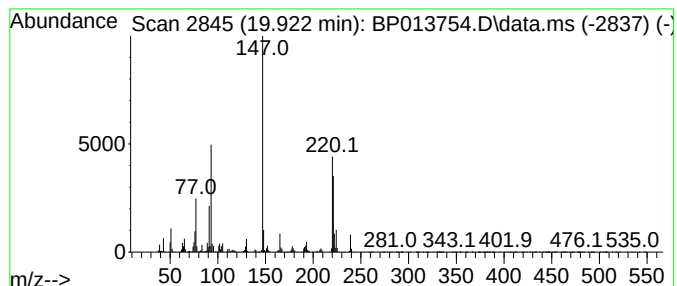
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 unknown-01 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.922	12.57 ng/ul	3206840	Chrysene-d12	21.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	3,4,8,8-Tetramethyl-4,5,6,7,8,8a...	220	C15H24O	1940169-10-7	43		
2	2H-Indazol-3-amine, 2-methyl-	147	C8H9N3	097990-19-7	43		
3	Ethanone, 1,1'-(1,3-phenylene)bis-	162	C10H10O2	006781-42-6	35		
4	Ethanone, 1,1'-(1,4-phenylene)bis-	162	C10H10O2	001009-61-6	35		
5	4-tert-Butylbenzyl mercaptan, S-...	222	C13H18OS	362528-98-1	27		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
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Sample : 01381-14
Misc :
ALS Vial : 55 Sample Multiplier: 1

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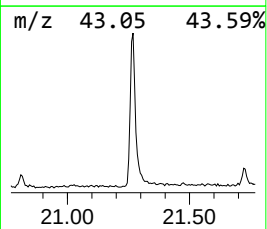
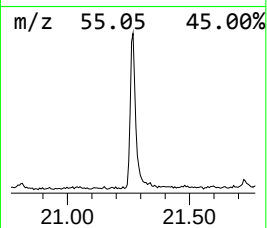
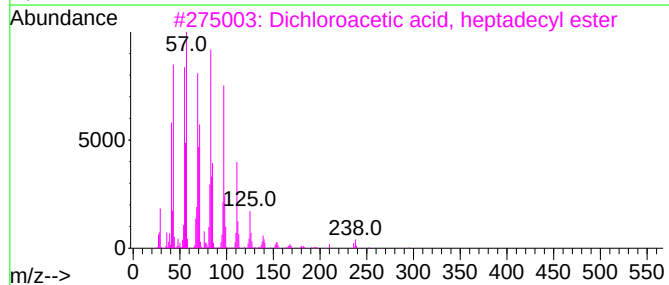
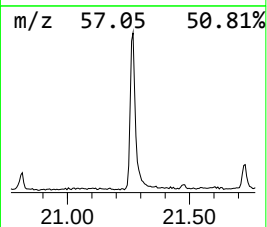
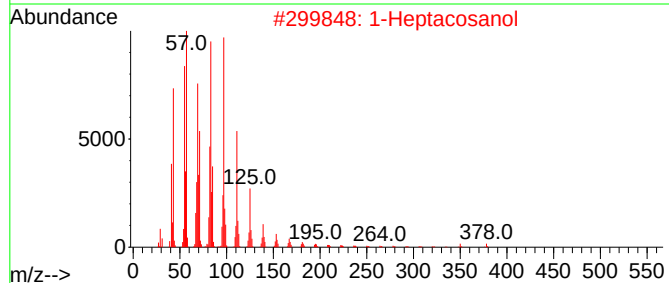
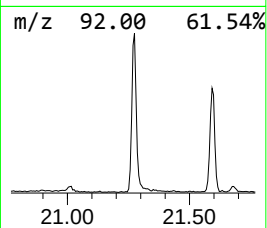
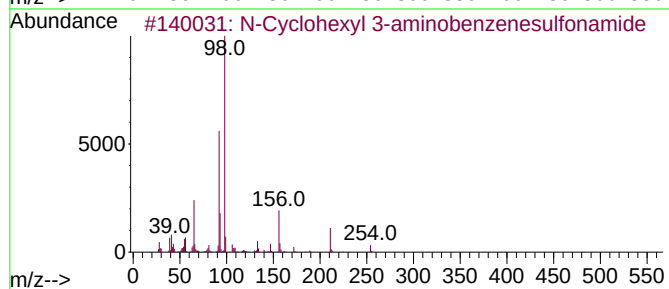
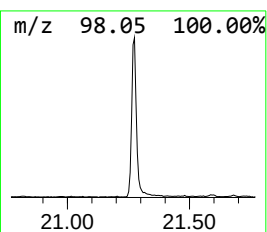
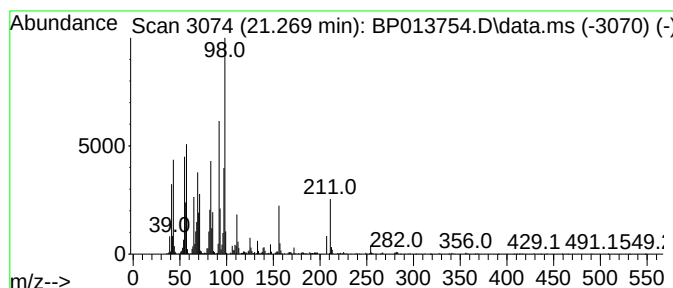
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 N-Cyclohexyl 3-aminobenzene... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.269	4.88 ng/ul	1245510	Chrysene-d12	21.598

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			N-Cyclohexyl 3-aminobenzenesulfo...	254	C12H18N2O2S	061886-26-8	95
2			1-Heptacosanol	396	C27H56O	002004-39-9	42
3			Dichloroacetic acid, heptadecyl ...	366	C19H36Cl2O2	1000282-98-2	42
4			Butyl tetracosyl ether	410	C28H58O	1000406-40-9	38
5			Butyl eicosyl ether	354	C24H50O	1000406-40-7	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
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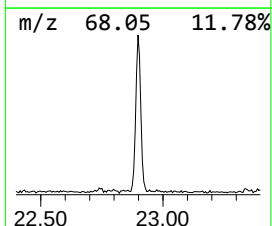
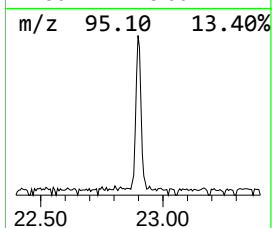
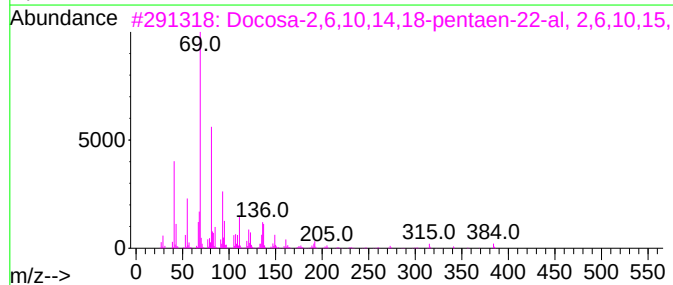
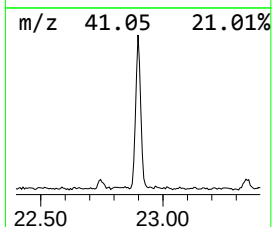
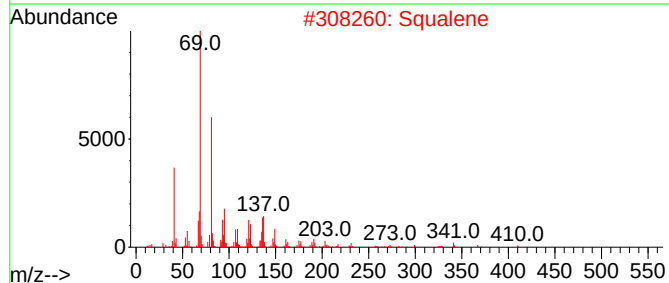
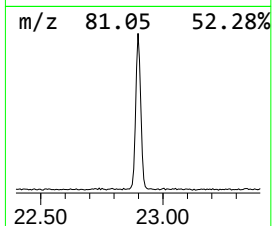
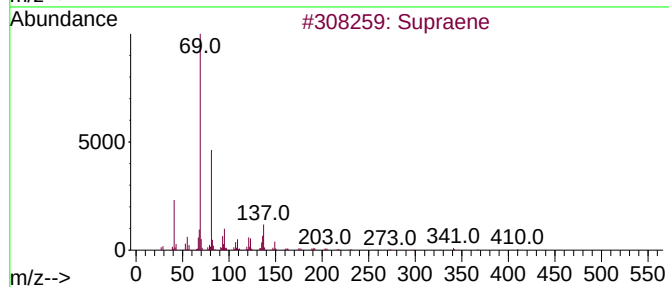
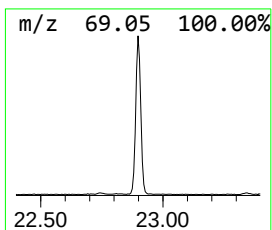
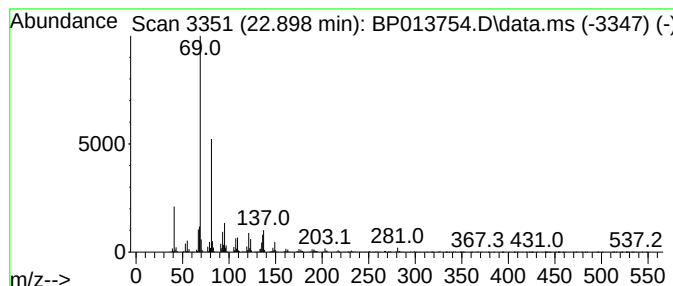
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Supraene Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.898	3.28 ng/ul	876344	Perylene-d12	24.163

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	98
2			Squalene	410	C30H50	000111-02-4	91
3			Docosa-2,6,10,14,18-pentaen-22-a...	384	C27H44O	1000163-04-7	90
4			Squalene	410	C30H50	000111-02-4	86
5			4-Methyl-1,5-Heptadiene	110	C8H14	000998-94-7	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP020223\
Data File : BP013754.D
Acq On : 03 Feb 2023 22:55
Operator : CG/JU
Sample : 01381-14
Misc :
ALS Vial : 55 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
A44N3

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP020123.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
Benzene, 1,3,5-...	10.816	17.0	ng/ul	1506420	2	10.957	1770370	20.0
Aniline, N-methyl-	12.463	3.3	ng/ul	289982	2	10.957	1770370	20.0
Benzophenone	16.116	10.4	ng/ul	1448940	3	14.763	2791140	20.0
n-Hexadecanoic ...	18.369	2.4	ng/ul	466535	4	17.516	3884030	20.0
(DEL) Alkane: S...	19.198	2.5	ng/ul	478938	4	17.516	3884030	20.0
5-Fluoroquinazo...	19.410	2.3	ng/ul	454335	4	17.516	3884030	20.0
7-Phenyl[1,2,4]...	19.445	12.6	ng/ul	2445440	4	17.516	3884030	20.0
unknown-01	19.922	12.6	ng/ul	3206840	5	21.598	5102560	20.0
N-Cyclohexyl 3-...	21.269	4.9	ng/ul	1245510	5	21.598	5102560	20.0
Supraene	22.898	3.3	ng/ul	876344	6	24.163	5349720	20.0