

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
 Data File : BP017524.D
 Acq On : 03 Oct 2023 19:14
 Operator : MA/JU
 Sample : 04679-26
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 PMW-7D(20230928)

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-BP092023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP017524.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.287	30	34	36	rBV	22915	28693	0.40%	0.062%
2	3.316	36	39	47	rVB	62471	79605	1.12%	0.171%
3	3.593	83	86	94	rVB	11288	16115	0.23%	0.035%
4	3.722	105	108	121	rBV	187839	267198	3.76%	0.575%
5	3.910	136	140	144	rVB5	8180	13220	0.19%	0.028%
6	4.028	156	160	165	rVB	14157	17580	0.25%	0.038%
7	4.446	228	231	235	rVB	16576	20603	0.29%	0.044%
8	4.552	244	249	254	rBV2	25699	33381	0.47%	0.072%
9	5.004	322	326	332	rBV	33246	45396	0.64%	0.098%
10	5.193	352	358	366	rVB	390184	561535	7.91%	1.208%
11	5.375	383	389	395	rBV	3755	7367	0.10%	0.016%
12	5.857	463	471	474	rBV6	7220	14375	0.20%	0.031%
13	6.622	597	601	610	rBV3	11497	18958	0.27%	0.041%
14	7.098	675	682	693	rBV	118078	204175	2.88%	0.439%
15	7.275	706	712	721	rBV	504610	756040	10.65%	1.626%
16	7.375	724	729	735	rVB2	11195	19920	0.28%	0.043%
17	7.457	735	743	752	rBV	499965	824155	11.61%	1.773%
18	7.534	752	756	764	rVB	28577	52136	0.73%	0.112%
19	7.810	796	803	815	rVB	962676	1445145	20.35%	3.108%
20	7.934	817	824	826	rBV	385481	594231	8.37%	1.278%
21	7.975	826	831	841	rVB	4580775	7099990	100.00%	15.270%
22	8.281	876	883	893	rVB	67337	112121	1.58%	0.241%
23	8.657	938	947	958	rBV	362654	614951	8.66%	1.323%
24	9.057	1010	1015	1019	rBV7	5047	8094	0.11%	0.017%
25	9.140	1019	1029	1037	rBV	594693	970965	13.68%	2.088%
26	9.263	1045	1050	1054	rVB8	4707	7100	0.10%	0.015%
27	9.340	1059	1063	1068	rVB5	12812	19265	0.27%	0.041%
28	9.463	1079	1084	1090	rBV2	25242	38976	0.55%	0.084%
29	9.522	1090	1094	1099	rVB5	8035	13793	0.19%	0.030%
30	9.857	1139	1151	1163	rBV	479216	802372	11.30%	1.726%
31	10.069	1183	1187	1193	rVB2	23282	36925	0.52%	0.079%
32	10.145	1195	1200	1207	rVB2	3478	7941	0.11%	0.017%
33	10.251	1214	1218	1223	rBV5	5077	8180	0.12%	0.018%
34	10.392	1234	1242	1251	rBV	692535	1197669	16.87%	2.576%
35	10.463	1251	1254	1262	rVB4	21835	38958	0.55%	0.084%
36	10.775	1300	1307	1314	rBV	502676	850015	11.97%	1.828%

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37	10.945	1325	1336	1357	rBV	522378	997525	14.05%	2.145%
38	11.422	1411	1417	1423	rBV3	7819	12208	0.17%	0.026%
39	11.545	1432	1438	1444	rBV3	4151	9781	0.14%	0.021%
40	12.116	1530	1535	1537	rBV6	4985	7818	0.11%	0.017%
41	12.298	1557	1566	1574	rBV	263377	446842	6.29%	0.961%
42	12.363	1574	1577	1584	rVB2	13735	24404	0.34%	0.052%
43	12.504	1594	1601	1606	rVB2	3641	7804	0.11%	0.017%
44	12.792	1646	1650	1658	rBV10	4718	12803	0.18%	0.028%
45	12.869	1658	1663	1668	rVB8	4679	8933	0.13%	0.019%
46	13.375	1746	1749	1752	rBV3	10623	12945	0.18%	0.028%
47	13.469	1759	1765	1773	rBV	63212	100371	1.41%	0.216%
48	13.616	1786	1790	1796	rVB8	5696	9637	0.14%	0.021%
49	13.686	1796	1802	1806	rBV9	6516	12443	0.18%	0.027%
50	13.810	1816	1823	1827	rBV4	10779	18373	0.26%	0.040%
51	13.986	1848	1853	1854	rBV5	5908	7633	0.11%	0.016%
52	14.045	1854	1863	1874	rVV	1431655	1915119	26.97%	4.119%
53	14.233	1891	1895	1903	rVB	29275	41550	0.59%	0.089%
54	14.328	1903	1911	1919	rBV2	1554599	2277204	32.07%	4.898%
55	14.416	1923	1926	1931	rVV2	15322	27808	0.39%	0.060%
56	14.451	1931	1932	1936	rVV4	7022	8441	0.12%	0.018%
57	14.522	1938	1944	1951	rVV	88270	122510	1.73%	0.263%
58	14.628	1954	1962	1970	rBV	702631	1041455	14.67%	2.240%
59	14.827	1992	1996	2000	rBV	17502	27459	0.39%	0.059%
60	14.898	2002	2008	2014	rBV2	48195	103803	1.46%	0.223%
61	15.022	2026	2029	2036	rVB8	11052	21209	0.30%	0.046%
62	15.210	2055	2061	2072	rBV	180802	281721	3.97%	0.606%
63	15.369	2083	2088	2093	rBV2	30718	47986	0.68%	0.103%
64	15.633	2126	2133	2140	rBV	2118734	2893045	40.75%	6.222%
65	15.704	2140	2145	2150	rVV	233200	324552	4.57%	0.698%
66	15.775	2152	2157	2169	rVV	385136	564787	7.95%	1.215%
67	15.875	2170	2174	2181	rVB10	8441	16469	0.23%	0.035%
68	16.180	2223	2226	2232	rBV3	12655	16564	0.23%	0.036%
69	16.298	2243	2246	2252	rVB8	5392	10780	0.15%	0.023%
70	16.769	2322	2326	2330	rBV6	14406	20159	0.28%	0.043%
71	17.333	2419	2422	2429	rVB6	15140	21924	0.31%	0.047%
72	17.421	2431	2437	2448	rBV	856702	1247076	17.56%	2.682%
73	17.521	2448	2454	2462	rBV	2458405	3314926	46.69%	7.129%
74	17.916	2517	2521	2524	rBV3	9741	13715	0.19%	0.029%
75	18.045	2539	2543	2548	rVB	20574	24078	0.34%	0.052%
76	18.145	2556	2560	2565	rBV8	18248	29771	0.42%	0.064%

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77	18.268	2576	2581	2594	rBV	506662	799735	11.26%	1.720%
78	19.186	2734	2737	2743	rBV5	27006	32082	0.45%	0.069%
79	19.345	2760	2764	2767	rBV2	29284	36551	0.51%	0.079%
80	19.398	2769	2773	2782	rVV3	81976	165969	2.34%	0.357%
81	19.515	2788	2793	2806	rBV2	150876	264042	3.72%	0.568%
82	19.774	2831	2837	2844	rBV	3203671	4030328	56.77%	8.668%
83	20.757	3002	3004	3010	rVB2	54347	64862	0.91%	0.139%
84	21.210	3076	3081	3090	rBV	463477	683238	9.62%	1.469%
85	21.557	3135	3140	3145	rBV	1093347	1412000	19.89%	3.037%
86	22.809	3349	3353	3361	rVB	329255	461619	6.50%	0.993%
87	23.968	3542	3550	3559	rBV2	1968117	4125919	58.11%	8.874%
88	24.139	3573	3579	3587	rVB	712324	1469242	20.69%	3.160%

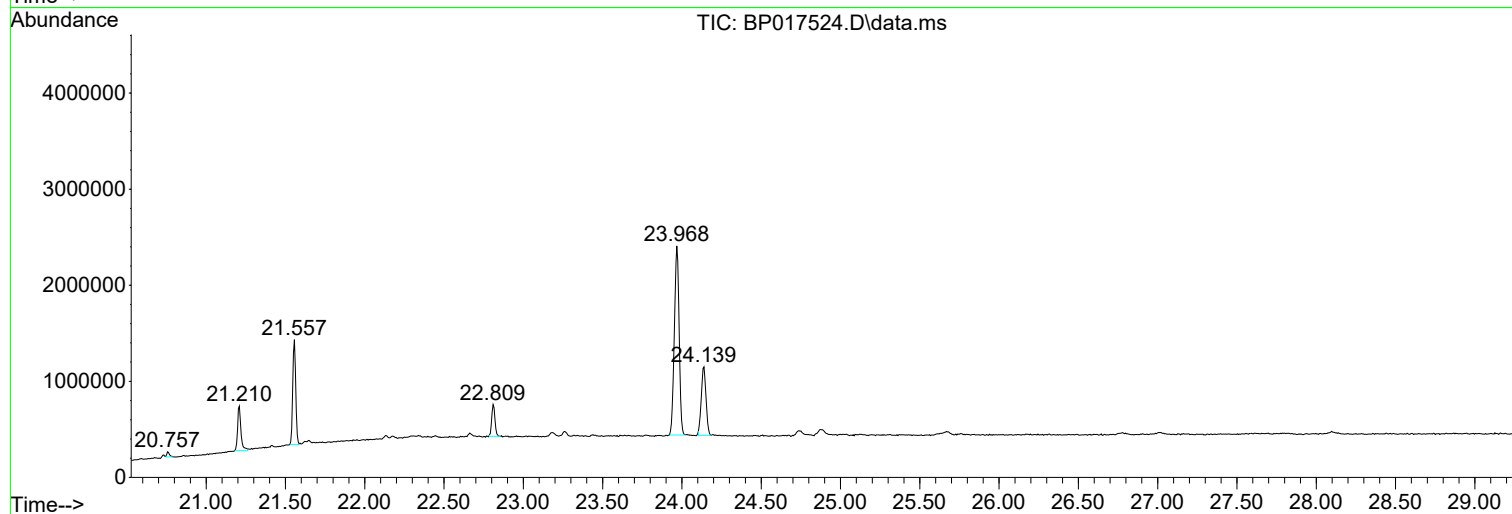
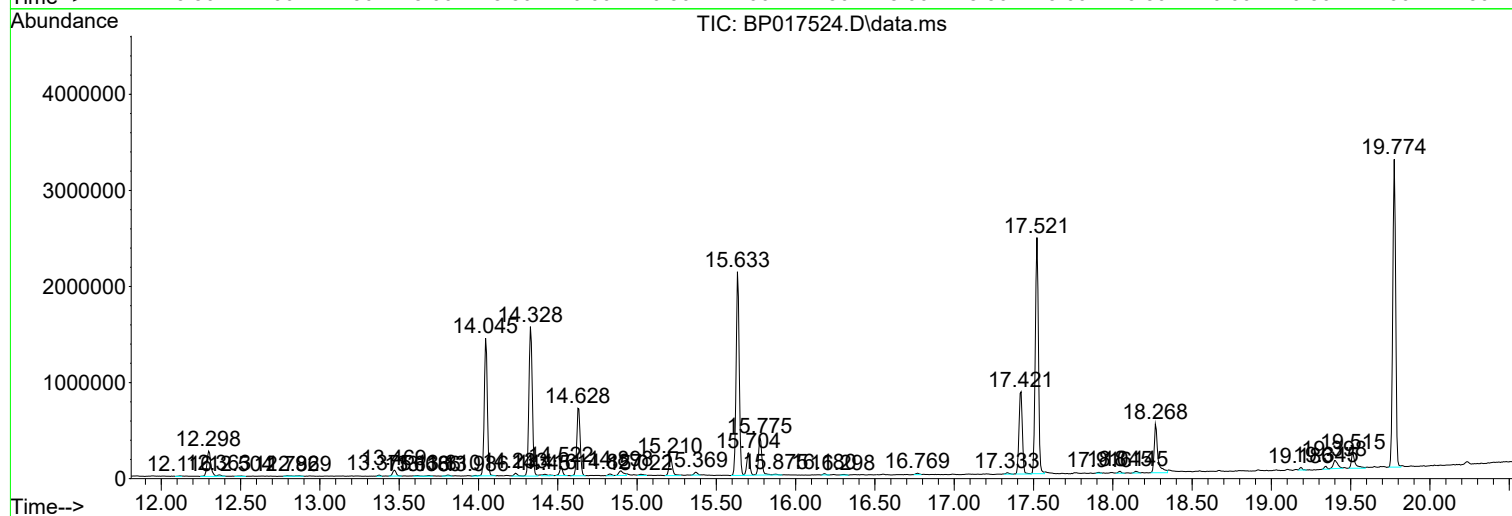
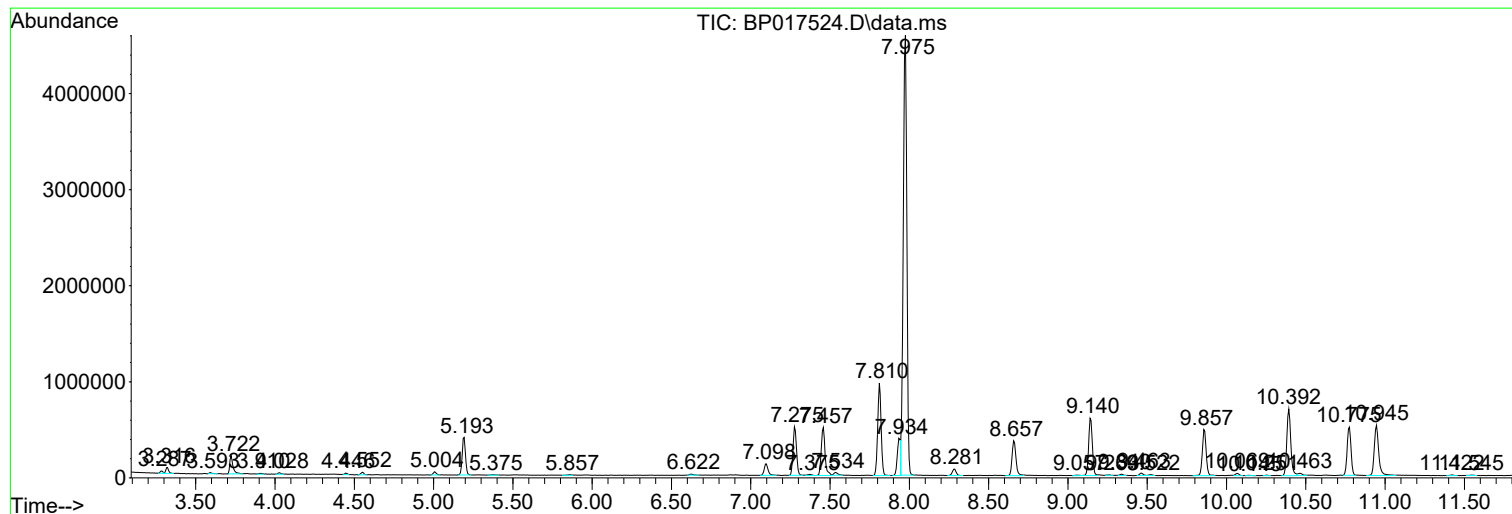
Sum of corrected areas: 46496361

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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



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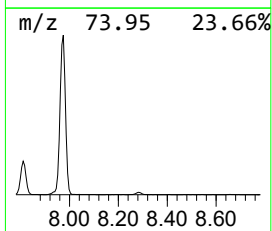
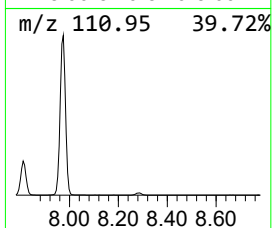
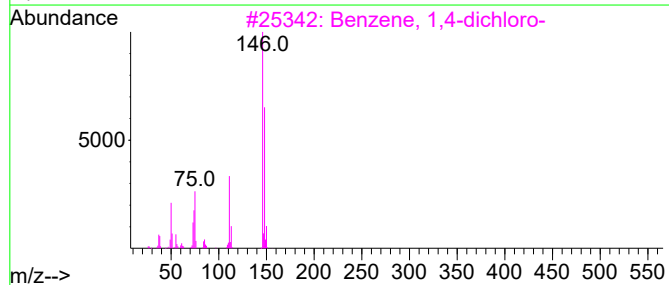
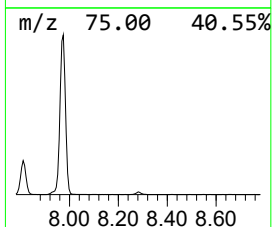
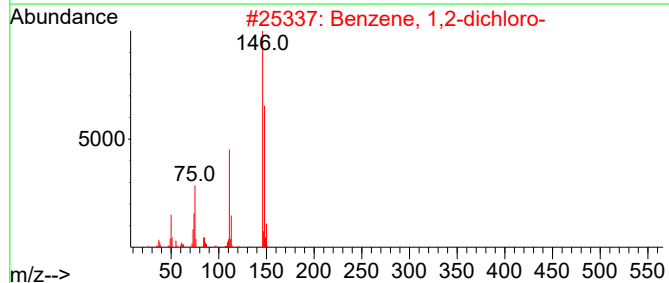
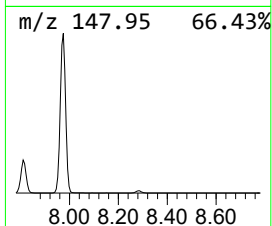
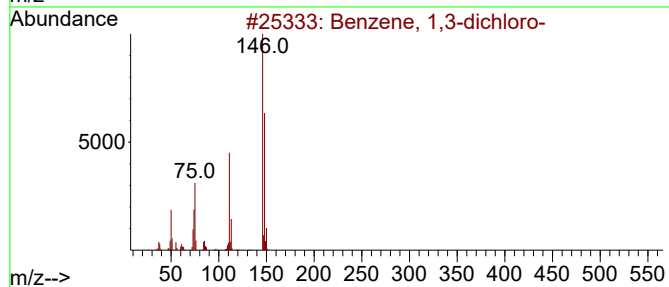
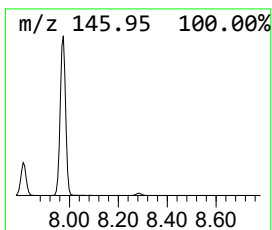
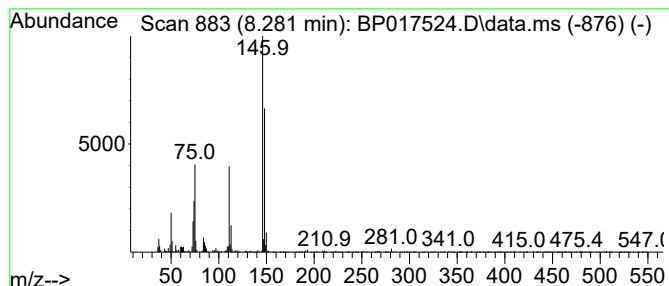
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, 1,3-dichloro- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.281	3.77 ng/ul	112121	1,4-Dichlorobenzene-d4	7.934

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	96
2			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	95
3			Benzene, 1,4-dichloro-	146	C6H4Cl2	000106-46-7	95
4			Benzene, 1,2-dichloro-	146	C6H4Cl2	000095-50-1	94
5			Benzene, 1,3-dichloro-	146	C6H4Cl2	000541-73-1	93



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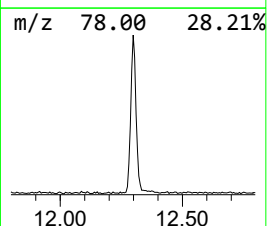
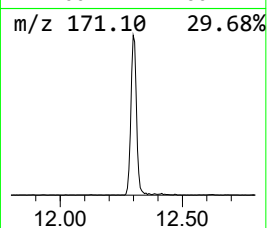
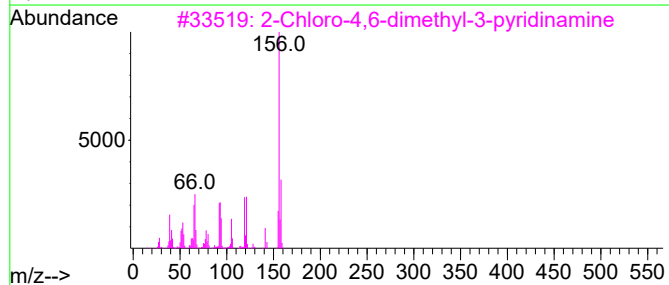
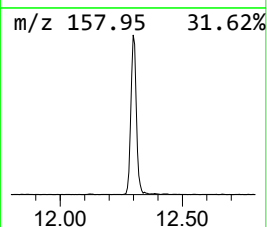
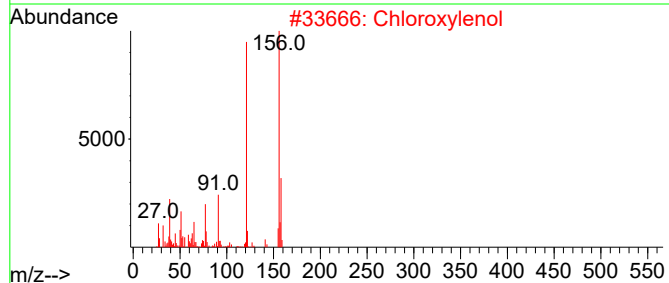
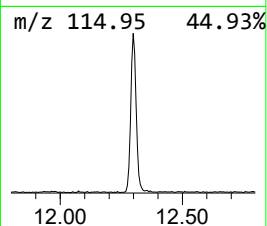
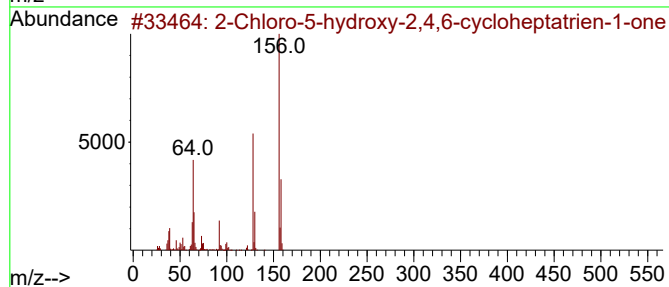
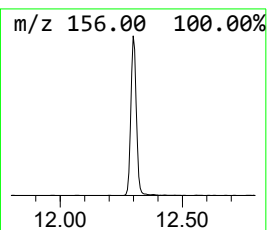
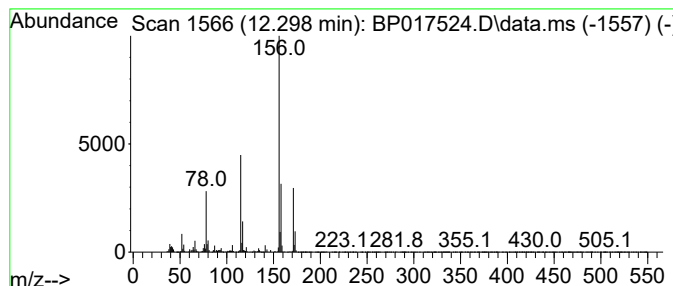
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 unknown-01 Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.298	10.51 ng/ul	446842	Naphthalene-d8	10.775

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Chloro-5-hydroxy-2,4,6-cyclohe...	156	C7H5ClO2	007009-16-7	43		
2	Chloroxylonol	156	C8H9ClO	000088-04-0	32		
3	2-Chloro-4,6-dimethyl-3-pyridina...	156	C7H9ClN2	140413-40-7	32		
4	Benzene, 2,5-cyclohexadien-1-yl-	156	C12H12	004794-05-2	25		
5	2-(2-Chloro-pyridin-3-yl)-propan...	171	C8H10ClNO	267003-35-0	25		



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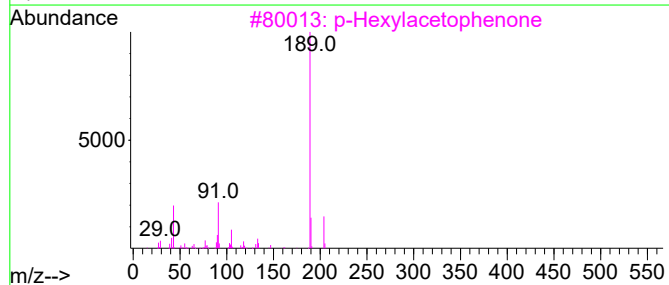
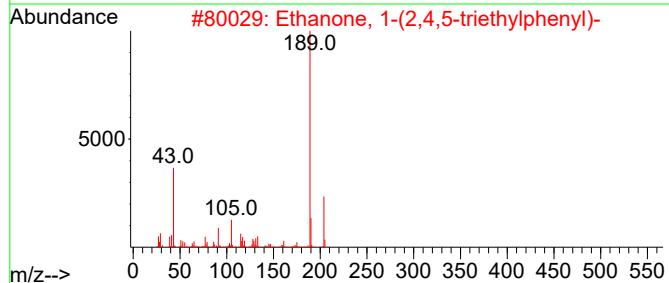
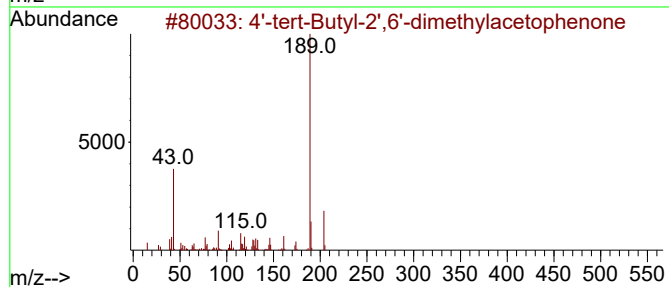
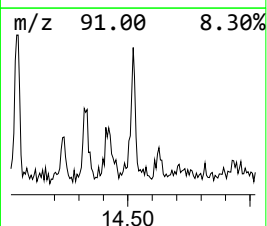
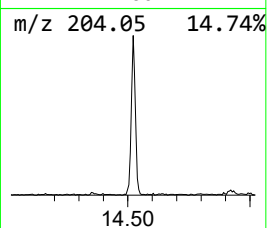
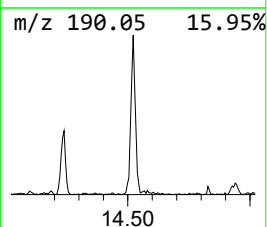
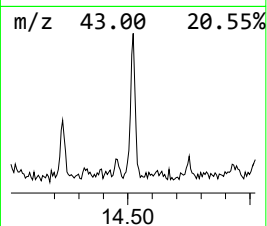
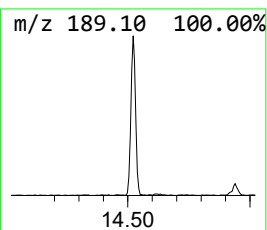
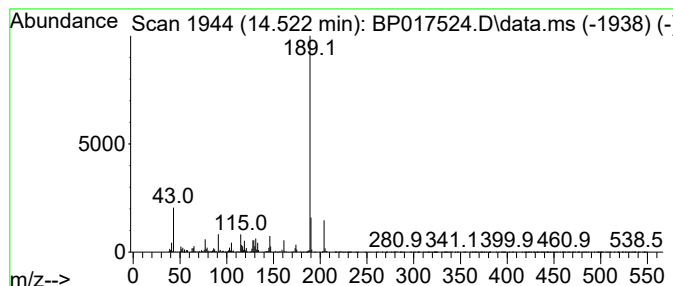
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4'-tert-Butyl-2',6'-dimethy... Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.522	2.35 ng/ul	122510	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4'-tert-Butyl-2',6'-dimethylacet...	204	C14H20O	002040-10-0	96
2			Ethanone, 1-(2,4,5-triethylphenyl)-	204	C14H20O	002715-54-0	90
3			p-Hexylacetophenone	204	C14H20O	037592-72-6	72
4			2-Nitro-1-naphthol	189	C10H7NO3	000607-24-9	64
5			Benzoic acid, 3-(3-hydroxy-3-met...	204	C12H12O3	1000337-06-2	64



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ALS Vial : 48 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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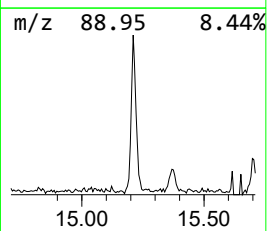
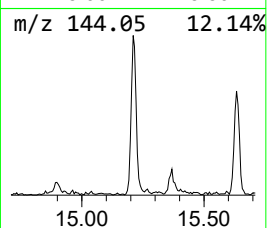
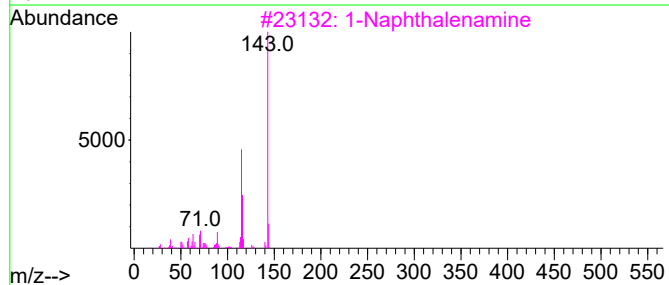
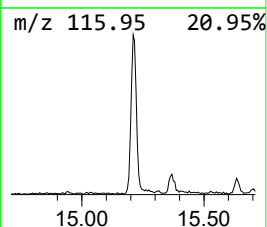
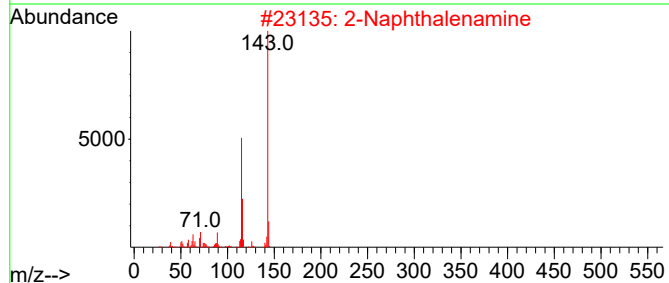
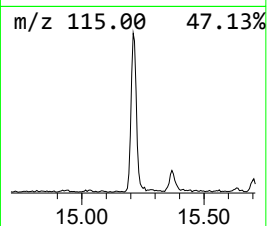
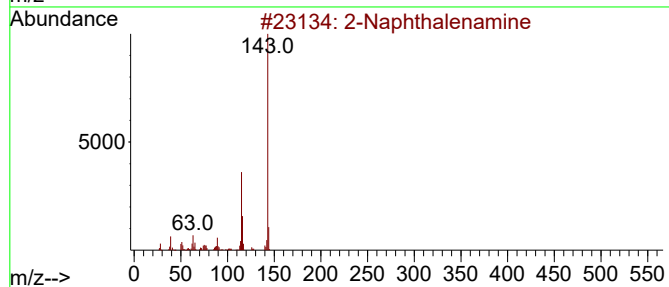
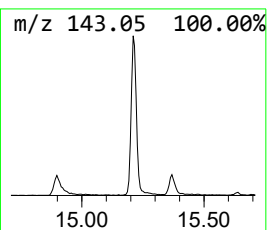
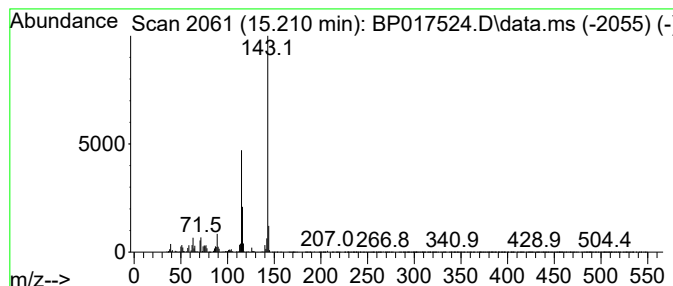
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Naphthalenamine Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.210	5.41 ng/ul	281721	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenamine			143	C10H9N	000091-59-8	97
2	2-Naphthalenamine			143	C10H9N	000091-59-8	97
3	1-Naphthalenamine			143	C10H9N	000134-32-7	97
4	1-Naphthalenamine			143	C10H9N	000134-32-7	96
5	2-Naphthalenamine			143	C10H9N	000091-59-8	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-7D(20230928)

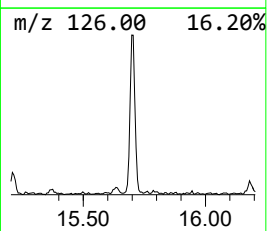
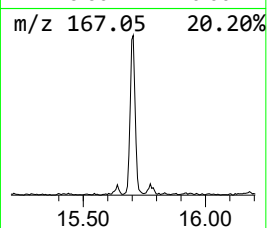
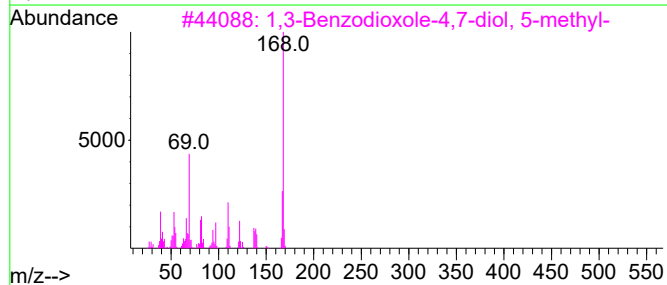
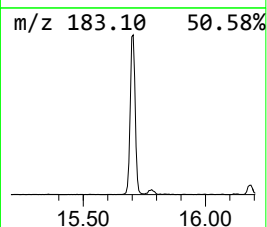
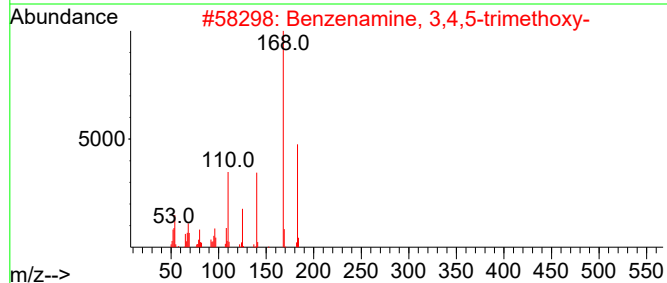
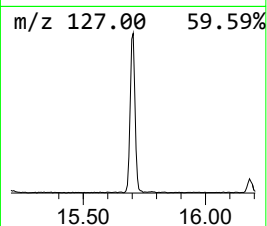
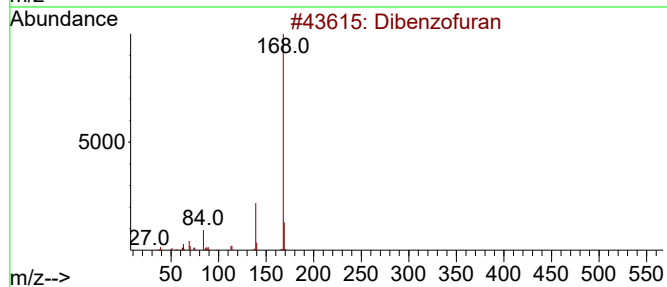
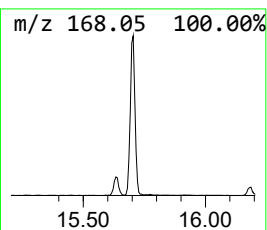
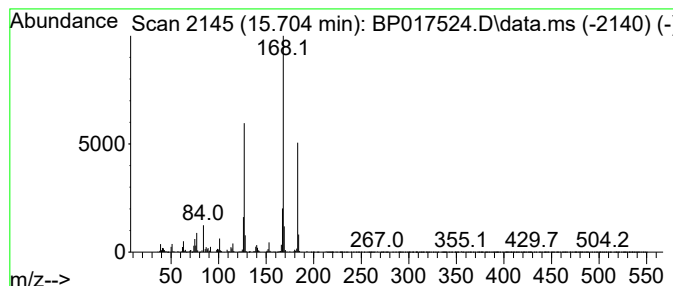
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 unknown-02 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.704	6.23 ng/ul	324552	Acenaphthene-d10	14.628

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Dibenzofuran	168	C12H8O	000132-64-9	46
2			Benzenamine, 3,4,5-trimethoxy-	183	C9H13NO3	024313-88-0	43
3			1,3-Benzodioxole-4,7-diol, 5-met...	168	C8H8O4	1187645-83-5	38
4			1-Naphthalenecarbonitrile, 8-amino-	168	C11H8N2	038515-13-8	35
5			Pyridine, 5-ethyl-2-phenyl-	183	C13H13N	066562-61-6	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-7D(20230928)

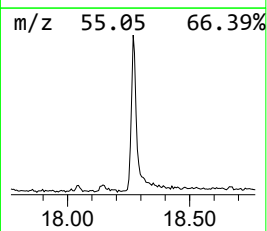
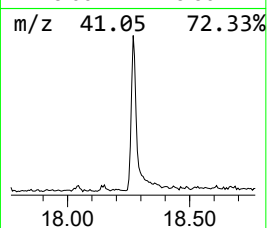
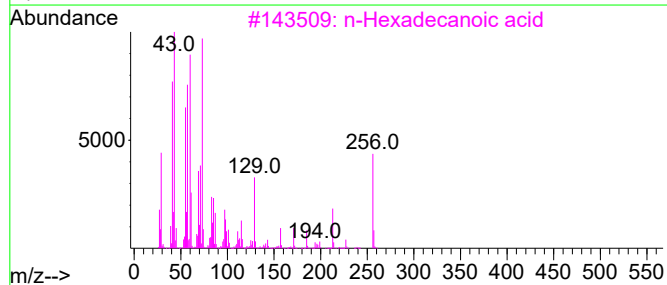
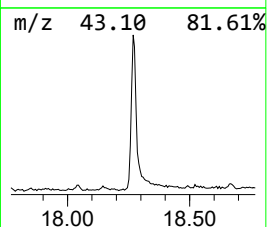
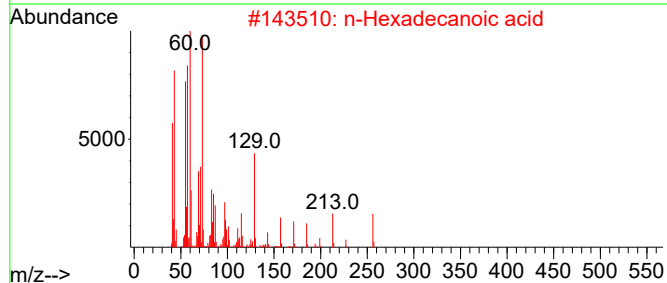
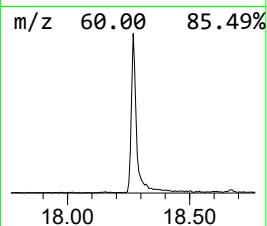
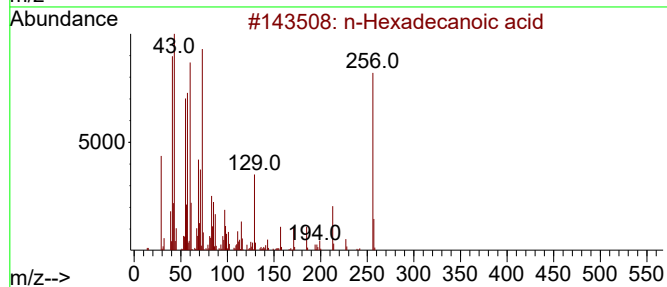
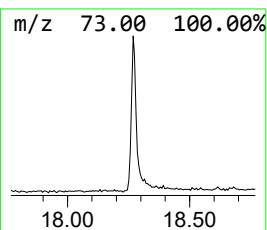
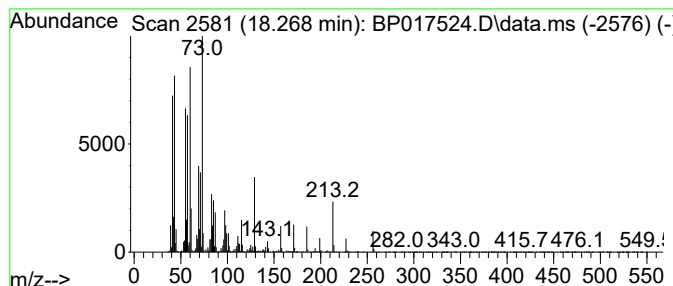
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 n-Hexadecanoic acid Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
18.269	12.83 ng/ul	799735	Phenanthrene-d10	17.422

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			Tridecanoic acid	214	C13H26O2	000638-53-9	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

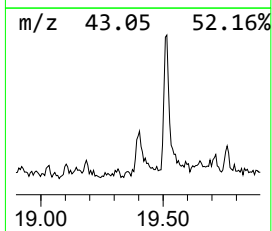
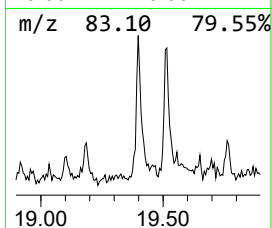
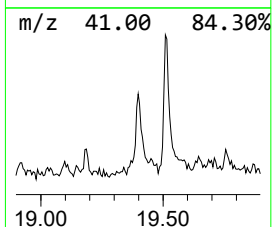
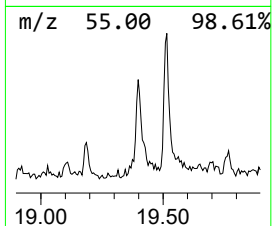
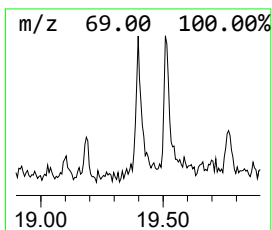
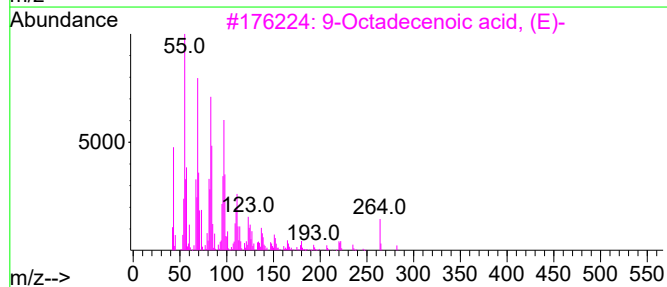
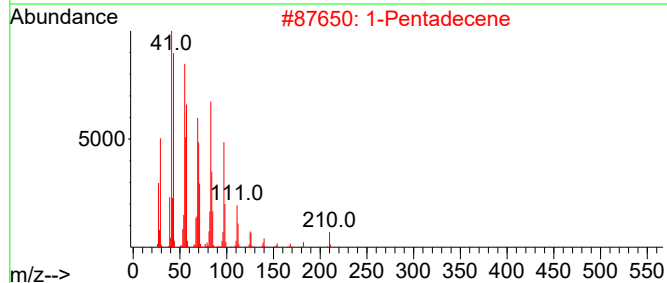
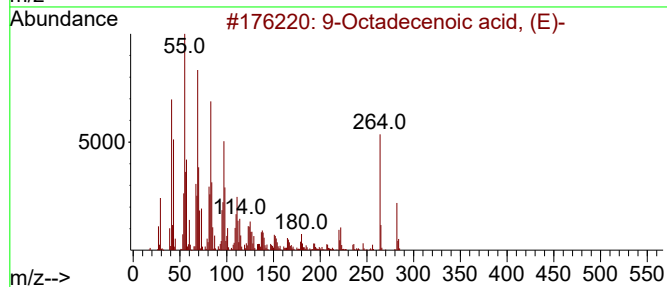
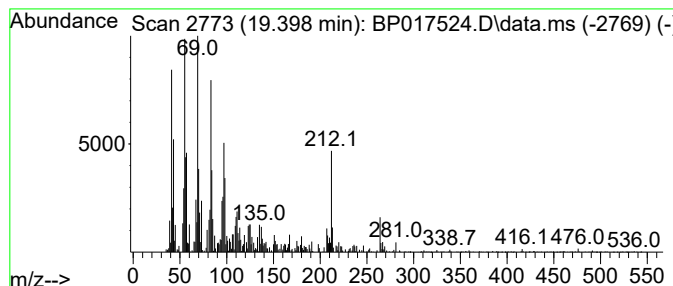
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 9-Octadecenoic acid, (E)- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.	
19.398	2.66 ng/ul	165969	Phenanthrene-d10	17.422	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	64
2	1-Pentadecene	210	C15H30	013360-61-7	60
3	9-Octadecenoic acid, (E)-	282	C18H34O2	000112-79-8	50
4	cis-9-Octadecenoic acid, propyl ...	324	C21H40O2	1000405-15-0	50
5	Cyclododecanol, 1-ethenyl-	210	C14H26O	006244-49-1	46



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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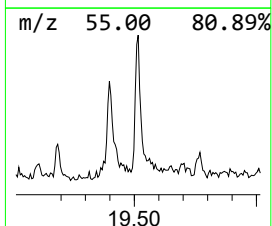
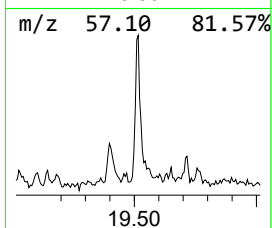
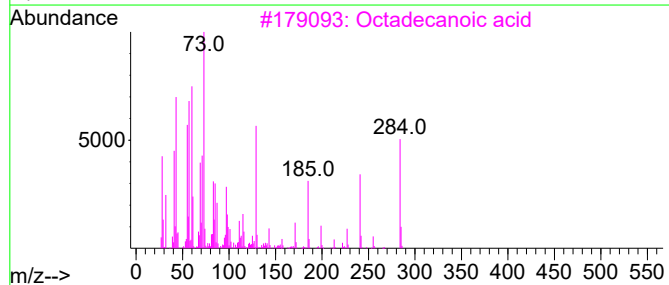
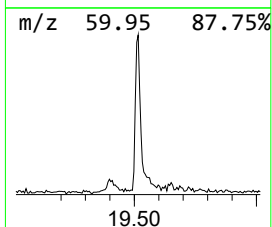
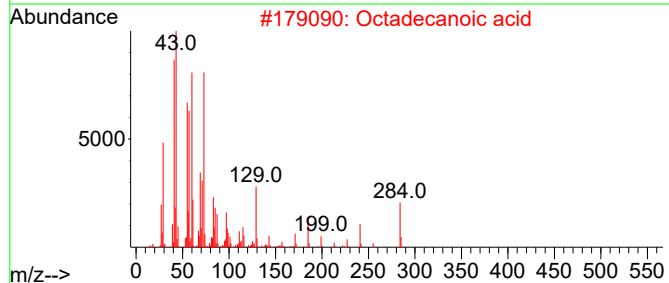
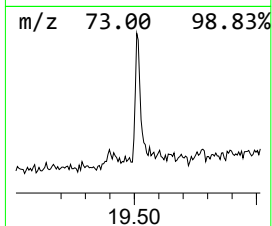
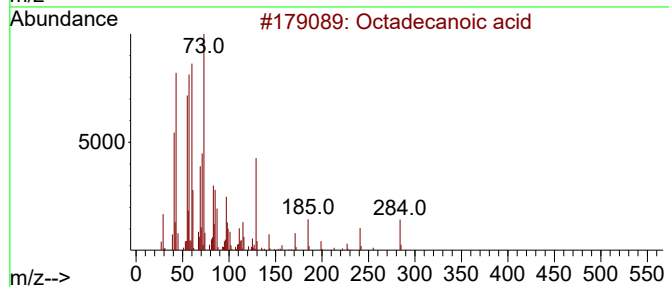
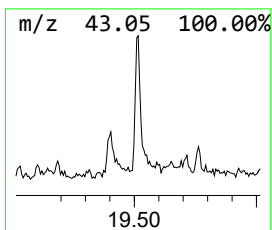
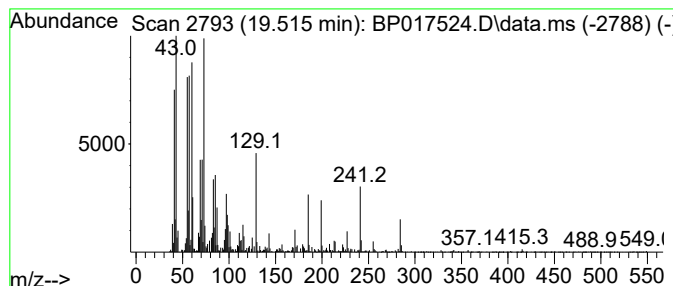
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Octadecanoic acid Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
19.515	3.74 ng/ul	264042	Chrysene-d12	21.557

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	97
3			Octadecanoic acid	284	C18H36O2	000057-11-4	97
4			Octadecanoic acid	284	C18H36O2	000057-11-4	96
5			n-Hexadecanoic acid	256	C16H32O2	000057-10-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
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ClientSampleId :
PMW-7D(20230928)

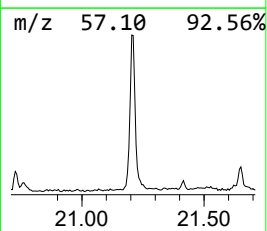
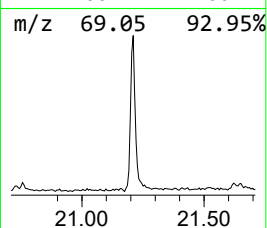
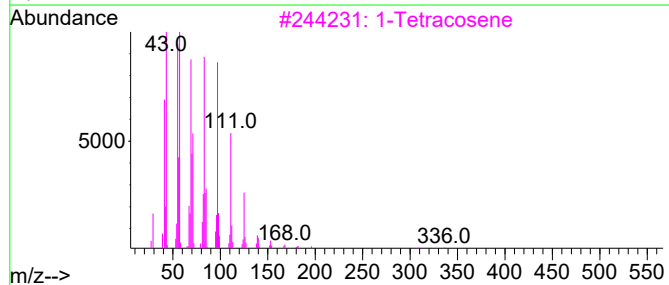
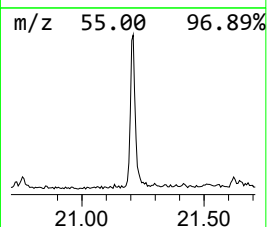
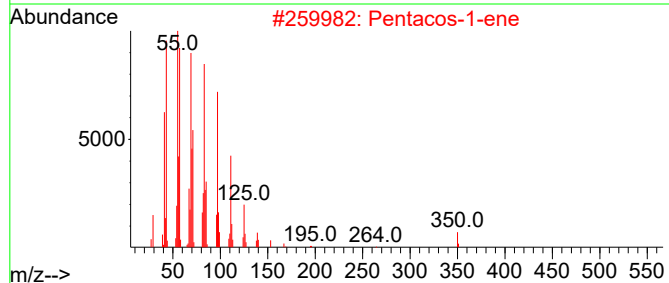
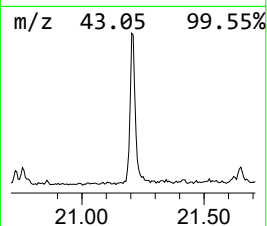
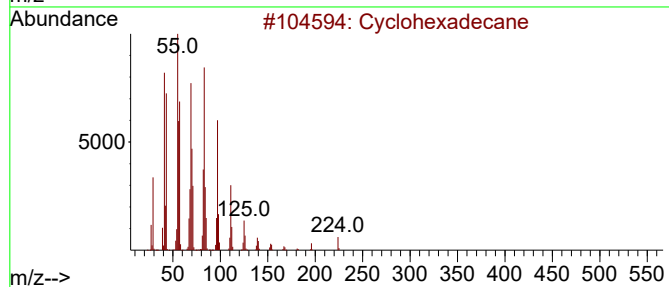
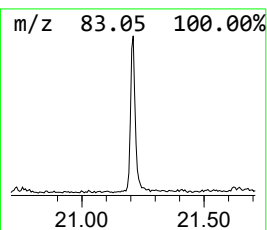
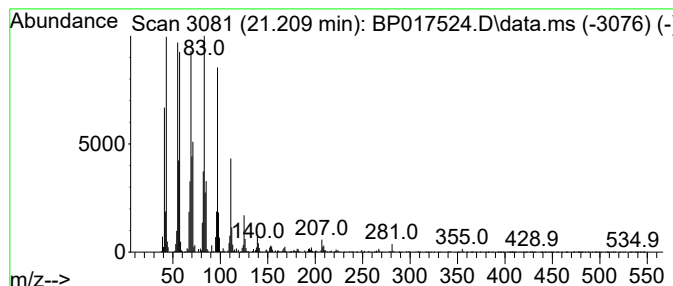
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 (DEL) Alkane: Cyclic21.209 Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
21.209	9.68 ng/ul	683238	Chrysene-d12	21.557

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		Cyclohexadecane	224	C16H32	000295-65-8	98
2		Pentacos-1-ene	350	C25H50	016980-85-1	94
3		1-Tetracosene	336	C24H48	010192-32-2	94
4		Heptafluorobutyric acid, pentade...	424	C19H31F7O2	959261-23-5	93
5		n-Nonadecanol-1	284	C19H40O	001454-84-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

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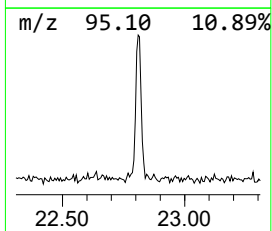
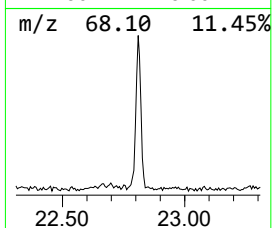
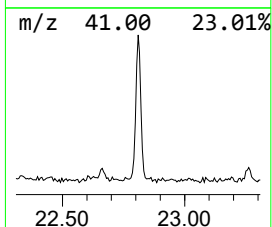
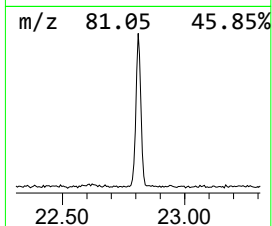
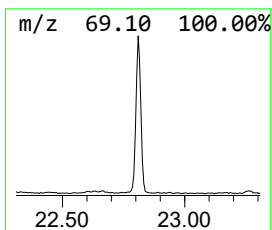
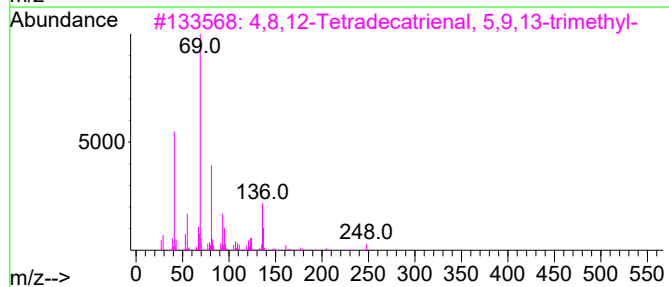
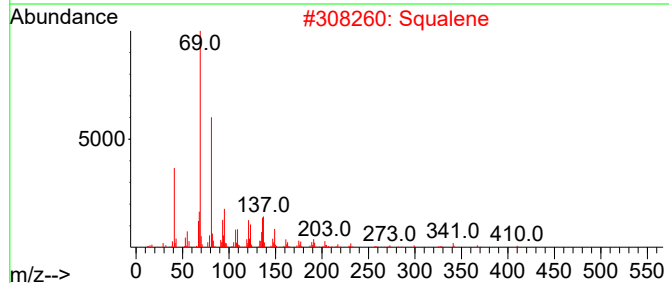
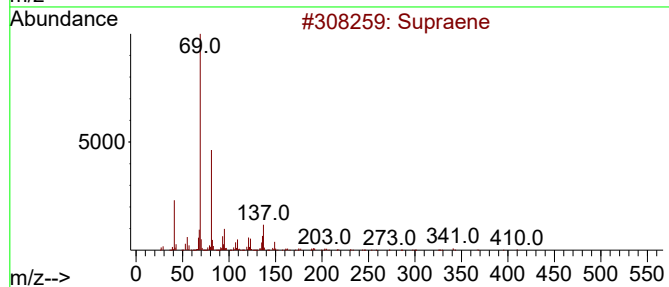
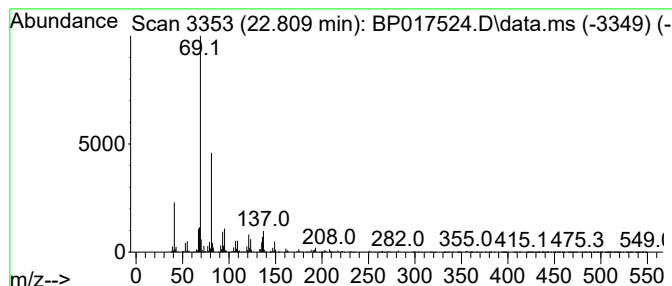
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 Supraene Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.809	6.54 ng/ul	461619	Chrysene-d12	21.557

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Supraene	410	C30H50	007683-64-9	91
2			Squalene	410	C30H50	000111-02-4	90
3			4,8,12-Tetradecatrienal, 5,9,13-...	248	C17H28O	066408-55-7	80
4			7,11-Dimethyldodeca-2,6,10-trien...	208	C14H24O	125529-10-4	72
5			1,5-Heptadiene, 3,3,6-trimethyl-	138	C10H18	035387-63-4	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP100223\
Data File : BP017524.D
Acq On : 03 Oct 2023 19:14
Operator : MA/JU
Sample : 04679-26
Misc :
ALS Vial : 48 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
PMW-7D(20230928)

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP092023.MA.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-di...	8.281	3.8	ng/ul	112121	1	7.934	594231	20.0
unknown-01	12.298	10.5	ng/ul	446842	2	10.775	850015	20.0
4'-tert-Butyl-2...	14.522	2.4	ng/ul	122510	3	14.628	1041460	20.0
2-Naphthalenamine	15.210	5.4	ng/ul	281721	3	14.628	1041460	20.0
unknown-02	15.704	6.2	ng/ul	324552	3	14.628	1041460	20.0
n-Hexadecanoic ...	18.269	12.8	ng/ul	799735	4	17.422	1247080	20.0
9-Octadecenoic ...	19.398	2.7	ng/ul	165969	4	17.422	1247080	20.0
Octadecanoic acid	19.515	3.7	ng/ul	264042	5	21.557	1412000	20.0
(DEL) Alkane: C...	21.209	9.7	ng/ul	683238	5	21.557	1412000	20.0
Supraene	22.809	6.5	ng/ul	461619	5	21.557	1412000	20.0