

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
 Data File : BM047014.D
 Acq On : 06 Aug 2024 06:52
 Operator : RC/JU
 Sample : P3382-01
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 E20R1

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_M\Methods\SFAM-EPA-BM080224.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BM047014.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	2.987	13	17	22	rBV	247614	320462	10.22%	1.042%
2	3.281	64	67	74	rBV	48007	70636	2.25%	0.230%
3	3.505	101	105	112	rVB	31884	42416	1.35%	0.138%
4	3.710	136	140	145	rVB	22955	32092	1.02%	0.104%
5	4.057	197	199	206	rVB9	10556	15585	0.50%	0.051%
6	4.469	262	269	273	rBV9	9155	15435	0.49%	0.050%
7	4.587	284	289	297	rBV3	27212	52492	1.67%	0.171%
8	5.693	471	477	482	rBV7	10452	17565	0.56%	0.057%
9	6.557	617	624	627	rBV	60275	99839	3.18%	0.325%
10	6.598	627	631	643	rVB	362785	592324	18.89%	1.927%
11	6.751	649	657	667	rBV	310275	460012	14.67%	1.496%
12	6.940	683	689	697	rVB	369765	576864	18.40%	1.876%
13	7.034	700	705	712	rBV3	29014	53977	1.72%	0.176%
14	7.393	759	766	775	rBV	1043790	1587059	50.61%	5.162%
15	7.475	775	780	788	rVB5	13097	26759	0.85%	0.087%
16	7.640	798	808	819	rBV	540009	904245	28.84%	2.941%
17	8.116	880	889	898	rBV	304158	486031	15.50%	1.581%
18	8.210	898	905	917	rVB	197346	322600	10.29%	1.049%
19	8.492	946	953	955	rBV6	8439	15049	0.48%	0.049%
20	8.540	955	961	968	rVV	353506	567032	18.08%	1.844%
21	8.616	971	974	979	rBV5	9402	13515	0.43%	0.044%
22	8.722	985	992	998	rVB5	8652	19117	0.61%	0.062%
23	9.122	1054	1060	1070	rVB	29597	51665	1.65%	0.168%
24	9.251	1072	1082	1090	rBV	261101	434370	13.85%	1.413%
25	9.439	1107	1114	1121	rBV	28264	54004	1.72%	0.176%
26	9.616	1139	1144	1149	rBV3	18186	29244	0.93%	0.095%
27	9.786	1166	1173	1181	rBV	409405	663255	21.15%	2.157%
28	10.028	1207	1214	1220	rVB8	9253	19654	0.63%	0.064%
29	10.151	1225	1235	1241	rBV	1298601	2091599	66.70%	6.803%
30	10.198	1241	1243	1249	rVB2	16076	22208	0.71%	0.072%
31	10.304	1252	1261	1274	rBV	261687	471107	15.02%	1.532%
32	10.728	1324	1333	1338	rBV8	18999	43863	1.40%	0.143%
33	10.910	1357	1364	1374	rBV	125296	241303	7.70%	0.785%
34	11.198	1408	1413	1417	rBV9	13371	24105	0.77%	0.078%
35	11.533	1464	1470	1475	rBV7	9037	14646	0.47%	0.048%
36	11.610	1478	1483	1491	rVV3	24367	41103	1.31%	0.134%

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Stop Thrs : 0

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37	11.751	1503	1507	1513	rVB4	15472	23675	0.75%	0.077%
38	11.828	1515	1520	1526	rBV3	18540	31738	1.01%	0.103%
39	12.051	1555	1558	1564	rVB2	12349	17042	0.54%	0.055%
40	12.457	1622	1627	1633	rVB8	11970	19378	0.62%	0.063%
41	12.539	1635	1641	1643	rBV7	8831	15176	0.48%	0.049%
42	12.592	1646	1650	1654	rVV	16922	23304	0.74%	0.076%
43	12.816	1683	1688	1694	rBV8	17845	47364	1.51%	0.154%
44	12.892	1696	1701	1707	rVV2	70568	112686	3.59%	0.367%
45	12.975	1711	1715	1719	rVV7	12429	20157	0.64%	0.066%
46	13.210	1750	1755	1760	rBV9	14066	33489	1.07%	0.109%
47	13.369	1777	1782	1786	rBV3	35541	52236	1.67%	0.170%
48	13.474	1794	1800	1807	rBV	778802	1112563	35.48%	3.619%
49	13.563	1812	1815	1818	rVV3	19827	24884	0.79%	0.081%
50	13.739	1838	1845	1850	rBV	867295	1322371	42.17%	4.301%
51	13.880	1865	1869	1874	rVB3	54401	67557	2.15%	0.220%
52	13.957	1878	1882	1883	rBV4	19177	24155	0.77%	0.079%
53	13.986	1883	1887	1890	rVV	94256	122800	3.92%	0.399%
54	14.051	1890	1898	1905	rVV	1863467	2715311	86.59%	8.832%
55	14.116	1905	1909	1911	rVV2	28772	44164	1.41%	0.144%
56	14.139	1911	1913	1919	rVB3	28769	47347	1.51%	0.154%
57	14.286	1931	1938	1939	rBV3	176278	225492	7.19%	0.733%
58	14.457	1960	1967	1972	rBV3	73128	154368	4.92%	0.502%
59	14.539	1979	1981	1985	rVB3	34530	41827	1.33%	0.136%
60	14.610	1990	1993	1997	rVB4	46468	56334	1.80%	0.183%
61	14.686	2003	2006	2009	rBV4	37384	44114	1.41%	0.143%
62	14.786	2019	2023	2029	rVB9	30665	50574	1.61%	0.164%
63	14.851	2029	2034	2039	rBV2	117750	182777	5.83%	0.594%
64	14.951	2047	2051	2054	rBV	99282	131289	4.19%	0.427%
65	15.057	2063	2069	2075	rBV	1203128	1686975	53.80%	5.487%
66	15.833	2195	2201	2206	rVB3	272213	504577	16.09%	1.641%
67	15.921	2211	2216	2222	rBV	541376	801903	25.57%	2.608%
68	15.980	2222	2226	2232	rVB2	459746	813655	25.95%	2.646%
69	16.045	2232	2237	2241	rBV2	453585	720976	22.99%	2.345%
70	16.086	2241	2244	2247	rVB	273992	278081	8.87%	0.904%
71	16.127	2247	2251	2255	rBV2	264366	427328	13.63%	1.390%
72	16.186	2259	2261	2265	rVB	217379	220585	7.03%	0.717%
73	16.239	2265	2270	2277	rBV3	454423	881113	28.10%	2.866%
74	16.310	2278	2282	2285	rBV	615986	802400	25.59%	2.610%
75	16.380	2291	2294	2300	rVB	371231	528820	16.86%	1.720%
76	16.810	2362	2367	2371	rBV	2201908	3135771	100.00%	10.199%

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Integrator: RTE
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Stop Thrs : 0 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
Peak separation: 5

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Title : SVOA CALIBRATION

77	16.851	2371	2374	2379	rVV	567275	749604	23.90%	2.438%
78	16.910	2379	2384	2394	rVV	1301234	1936177	61.74%	6.297%

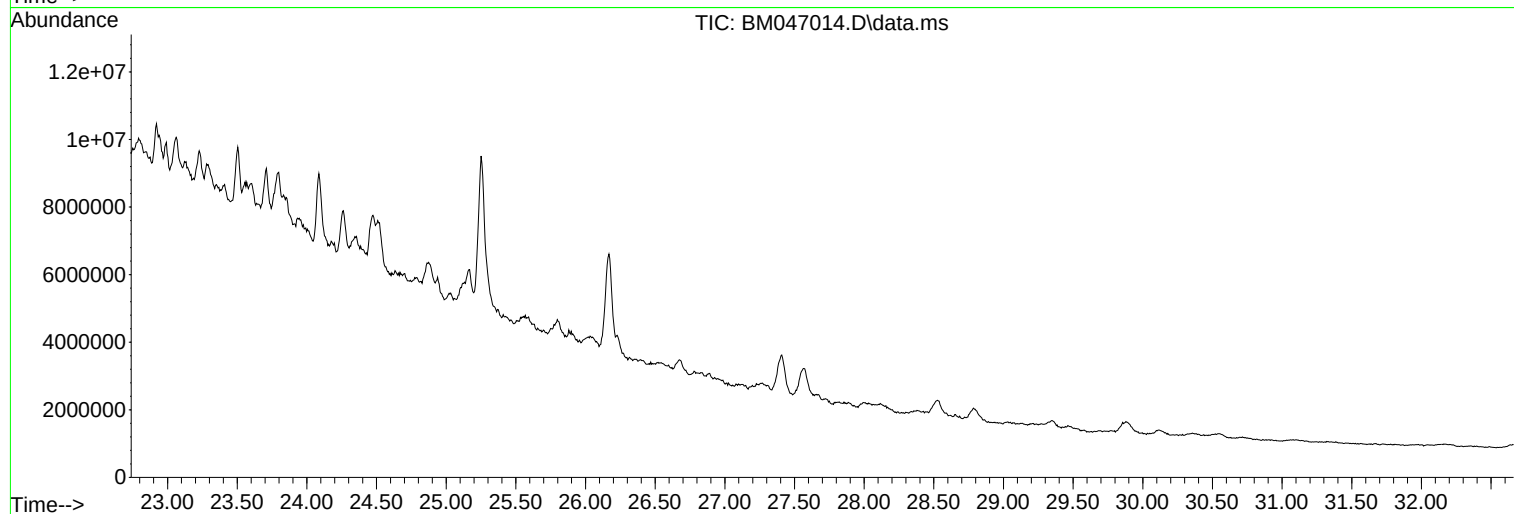
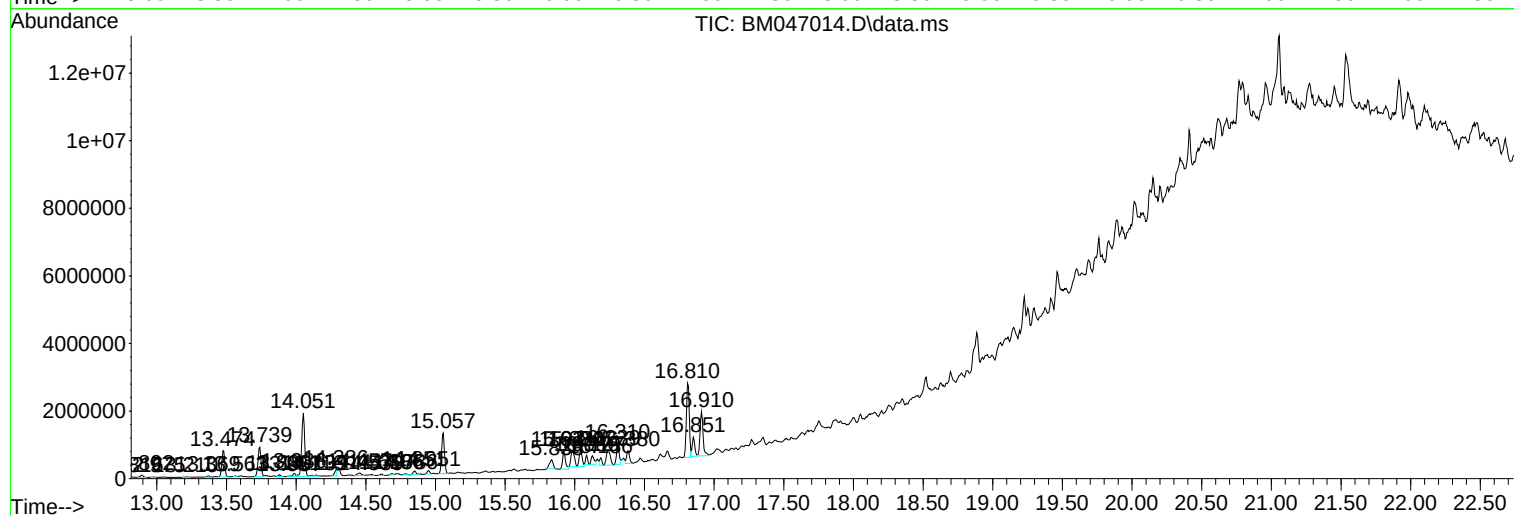
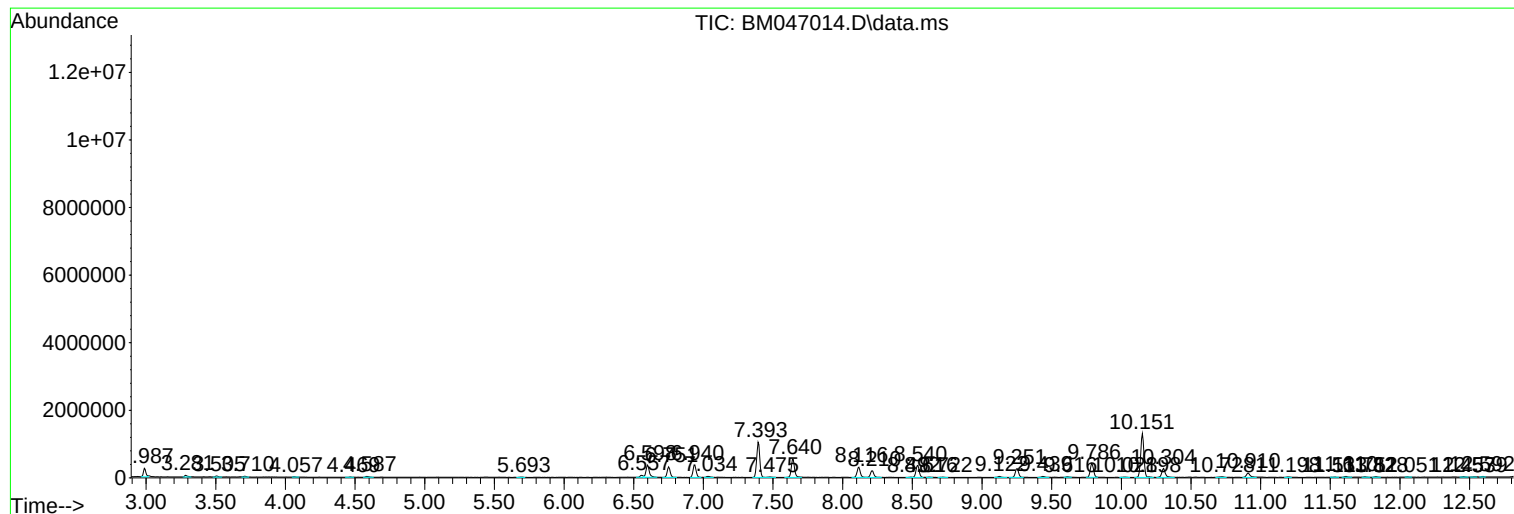
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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



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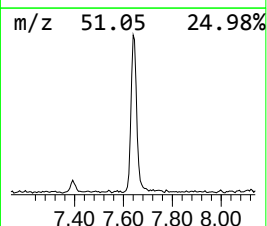
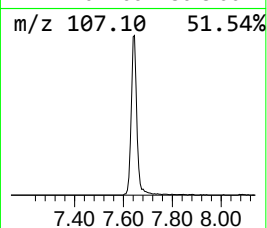
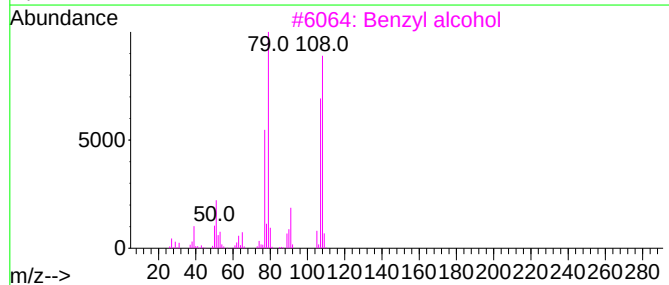
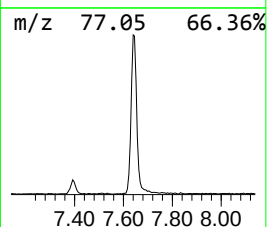
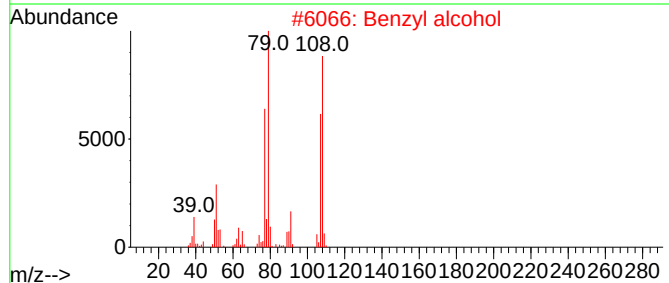
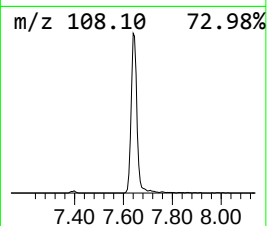
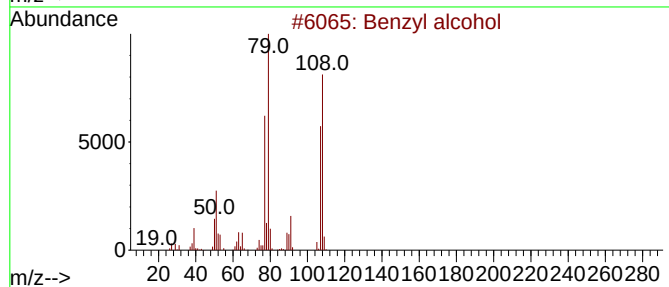
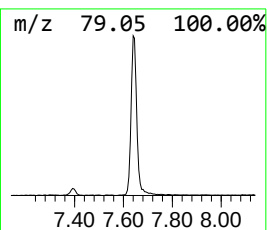
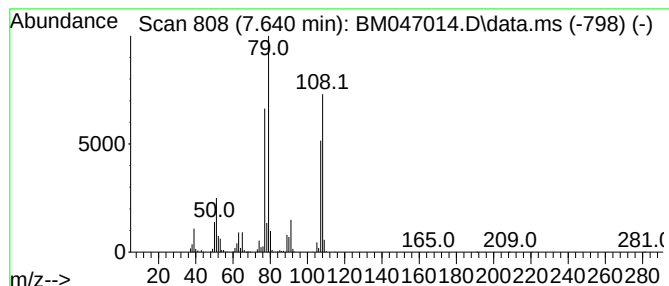
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzyl alcohol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.640	11.40 ng/ul	904245	1,4-Dichlorobenzene-d4	7.393

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzyl alcohol	108	C7H8O	000100-51-6	98
2			Benzyl alcohol	108	C7H8O	000100-51-6	98
3			Benzyl alcohol	108	C7H8O	000100-51-6	97
4			Benzyl alcohol	108	C7H8O	000100-51-6	96
5			N-Cbz-glycylglycine p-nitropheny...	387	C18H17N3O7	013574-81-7	91



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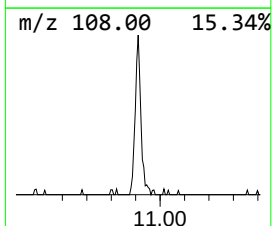
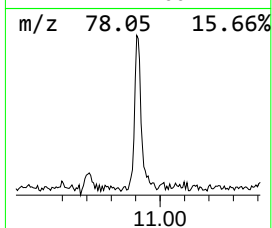
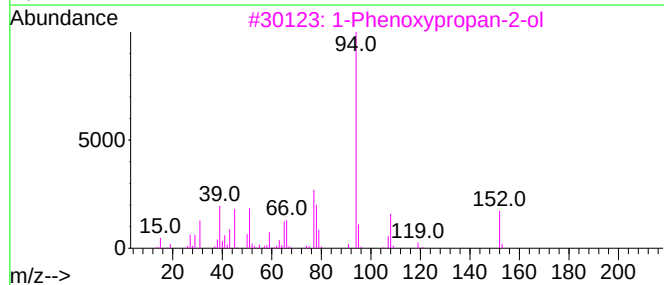
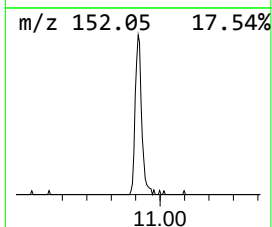
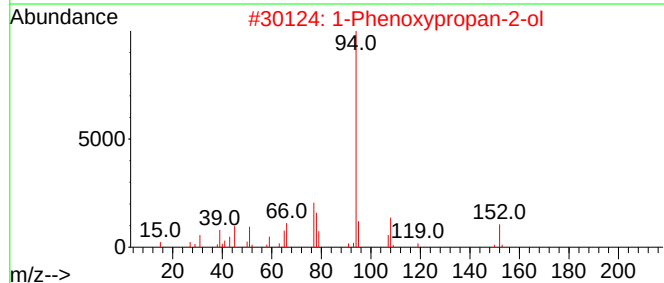
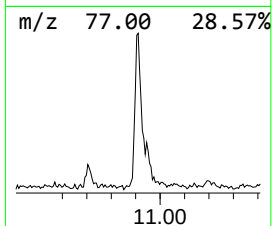
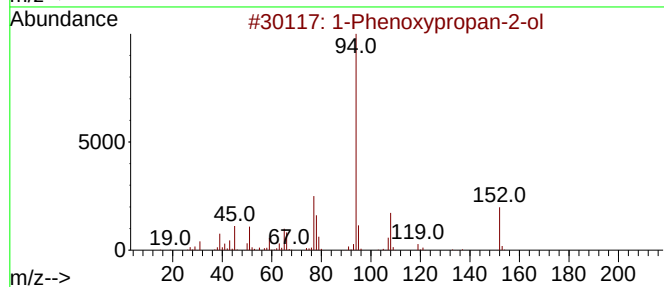
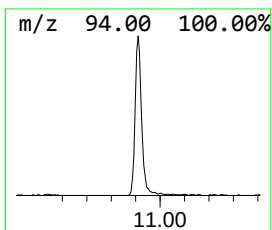
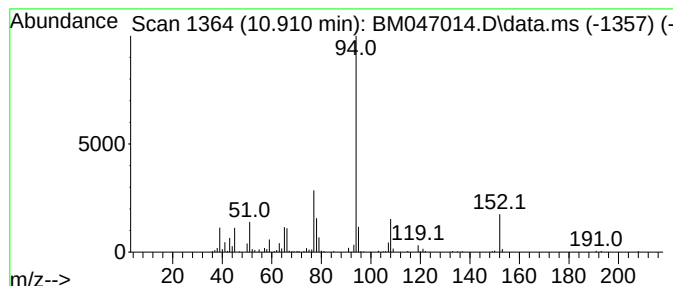
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Phenoxypropan-2-ol Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.910	2.31 ng/ul	241303	Naphthalene-d8	10.151

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	97
2			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
3			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	91
4			1-Phenoxypropan-2-ol	152	C9H12O2	000770-35-4	87
5			1-Propanol, 3-phenoxy-	152	C9H12O2	006180-61-6	83



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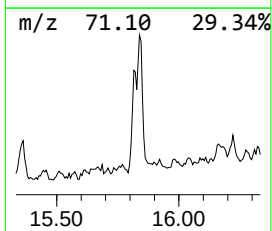
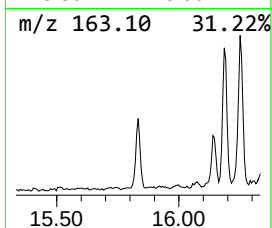
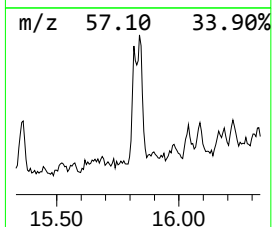
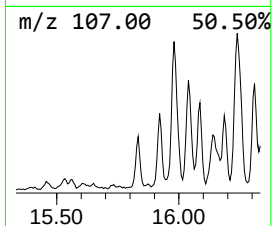
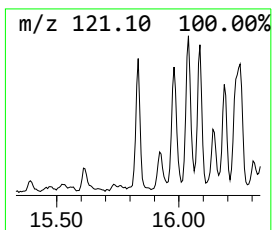
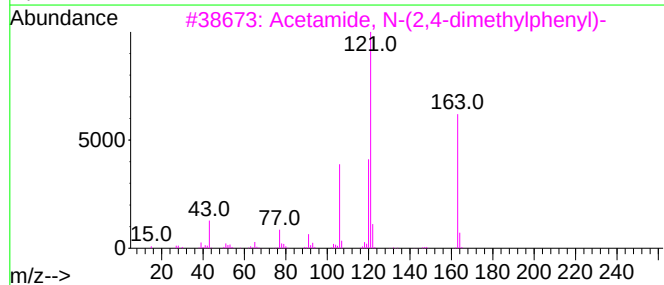
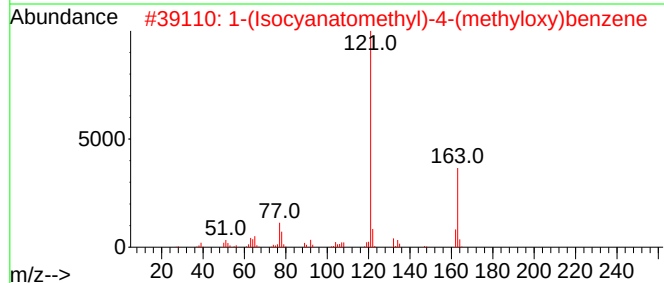
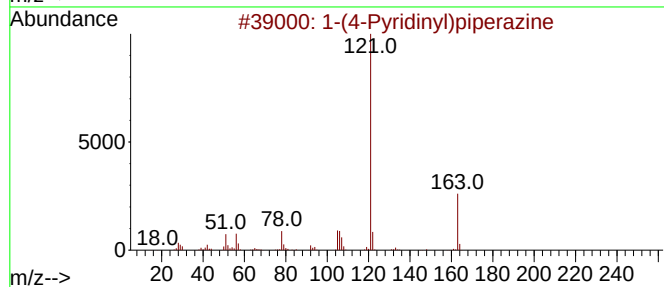
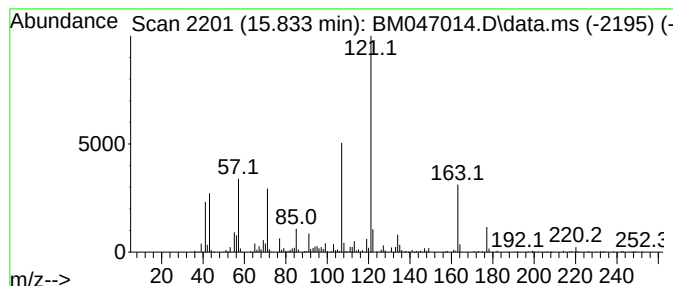
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 unknown-01 Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.833	3.22 ng/ul	504577	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(4-Pyridinyl)piperazine	163	C9H13N3	001008-91-9	49
2			1-(Isocyanatomethyl)-4-(methylox...	163	C9H9NO2	056651-60-6	47
3			Acetamide, N-(2,4-dimethylphenyl)-	163	C10H13NO	002050-43-3	46
4			Anisole, p-octyl-	220	C15H24O	003307-19-5	38
5			1-Chloro-1-cyclopropyl-2-(4-meth...	226	C12H15ClO2	1000486-82-8	38



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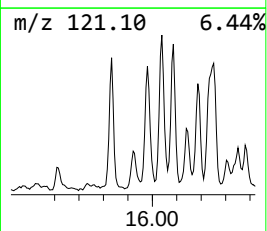
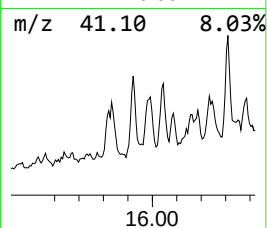
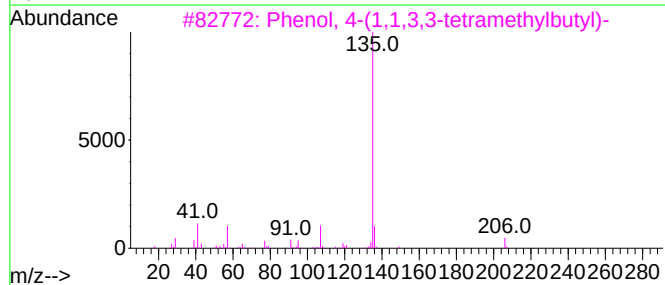
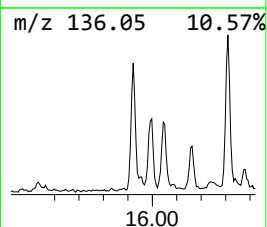
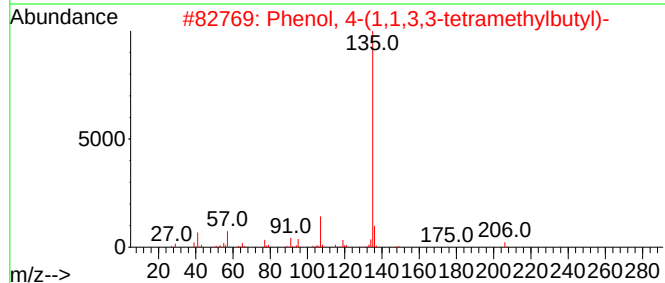
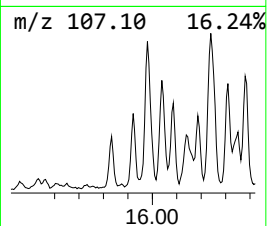
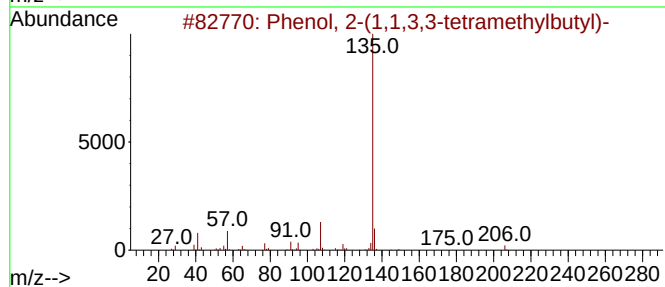
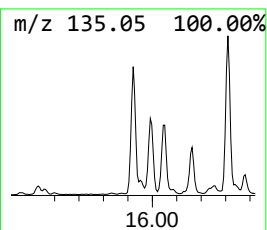
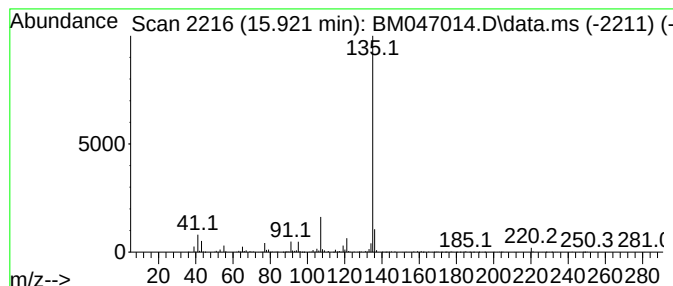
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Phenol, 2-(1,1,3,3-tetramet... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.921	5.11 ng/ul	801903	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	86
2			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
3			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
4			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
5			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
Operator : RC/JU
Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E20R1

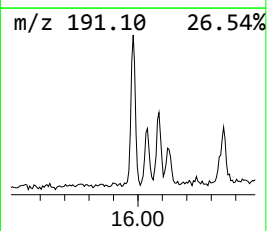
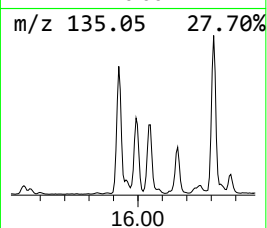
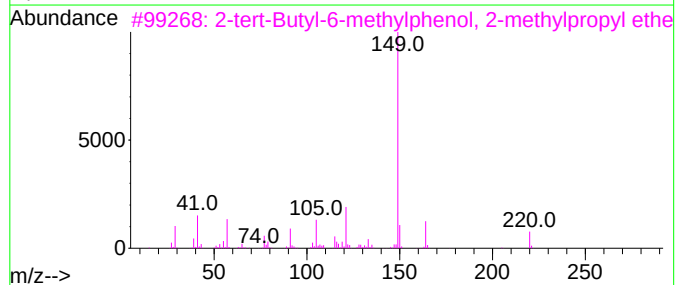
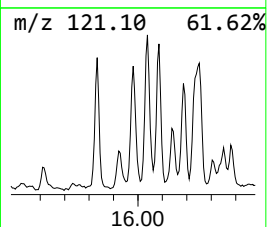
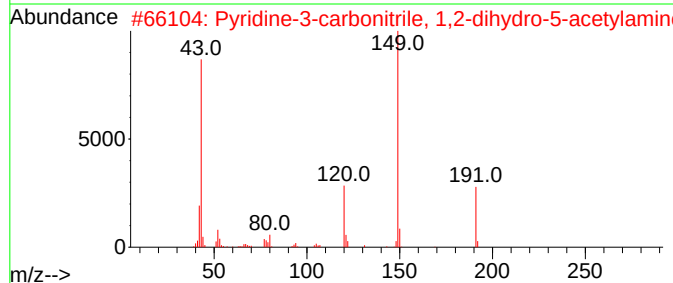
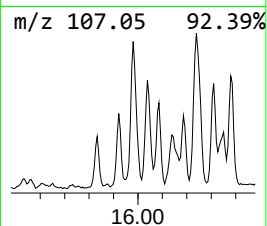
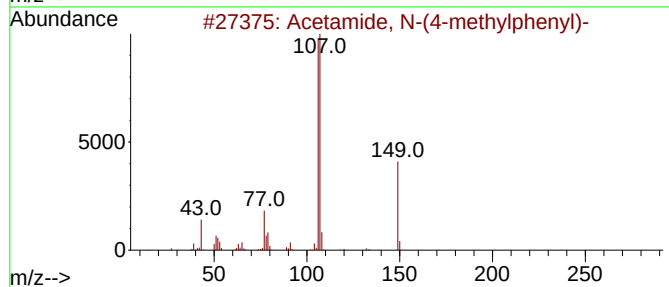
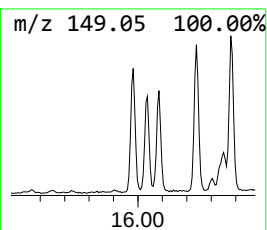
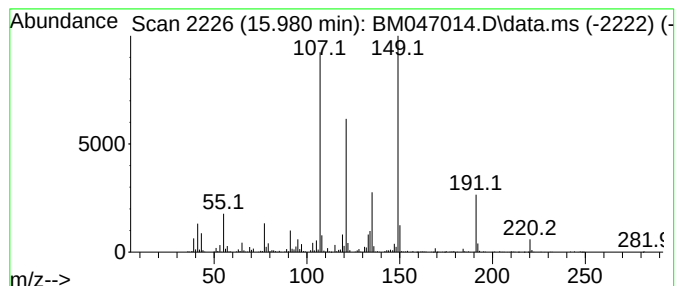
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 unknown-02 Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.980	5.19 ng/ul	813655	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(4-methylphenyl)-	149	C9H11NO	000103-89-9	43
2			Pyridine-3-carbonitrile, 1,2-dih...	191	C9H9N3O2	220098-92-0	38
3			2-tert-Butyl-6-methylphenol, 2-m...	220	C15H24O	1000395-19-4	35
4			Phthalic acid, phenyl propyl ester	284	C17H16O4	1000314-86-9	35
5			Benzene, 1-butyl-4-methoxy-	164	C11H16O	018272-84-9	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
Operator : RC/JU
Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E20R1

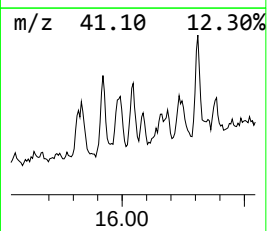
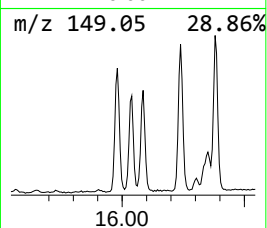
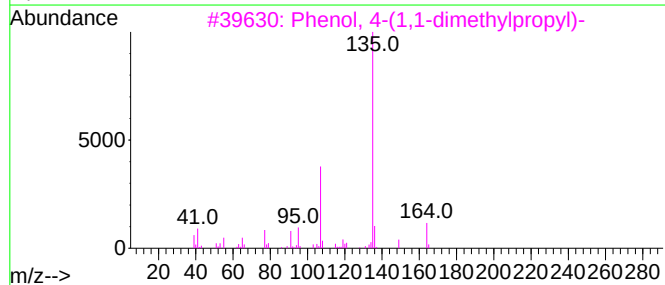
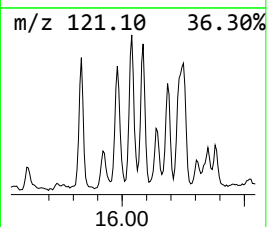
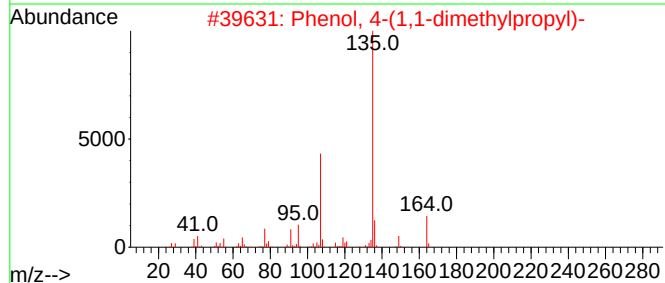
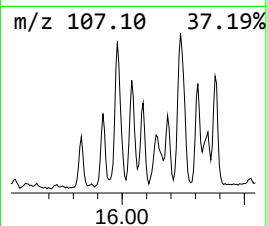
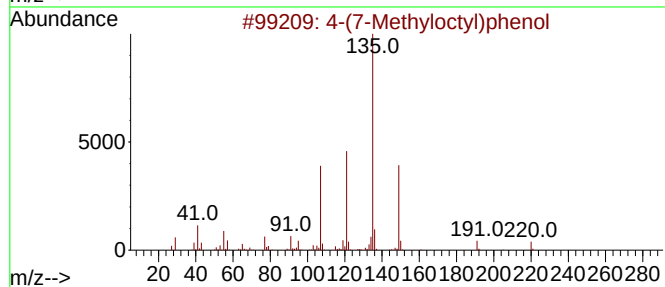
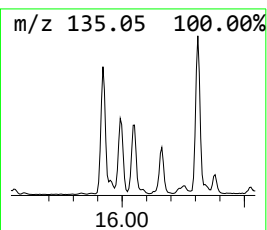
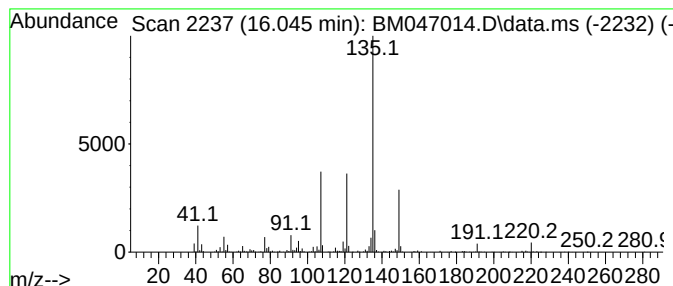
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 4-(7-Methyloctyl)phenol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.045	4.60 ng/ul	720976	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	95
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	64
4			Hexestrol, 0-methoxycarbonyl-	328	C20H24O4	1000487-66-3	59
5			4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	59



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
Operator : RC/JU
Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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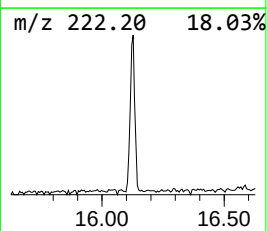
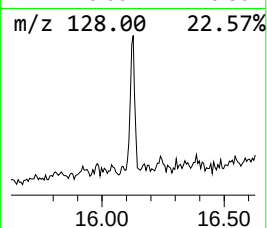
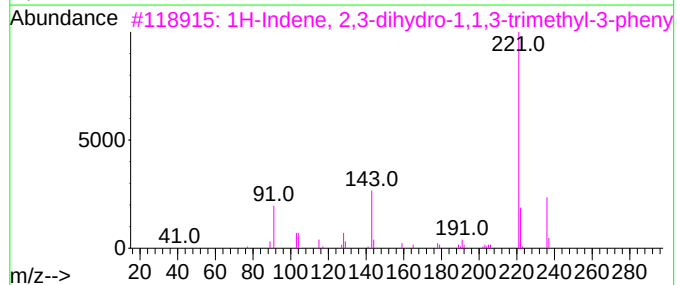
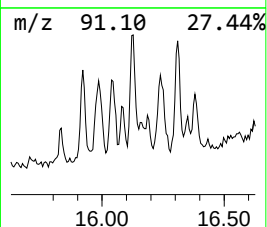
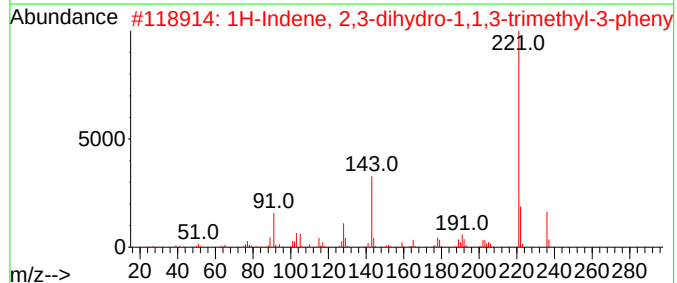
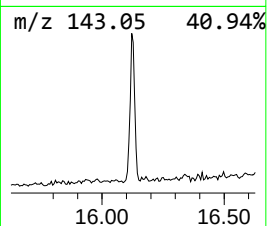
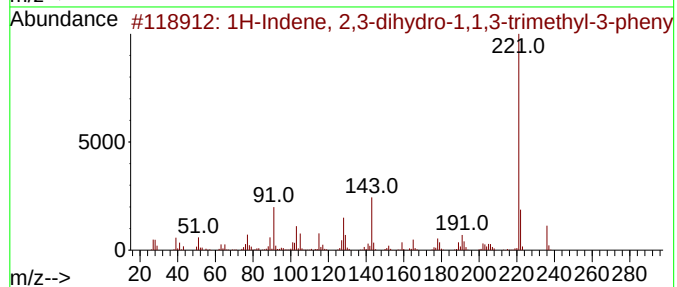
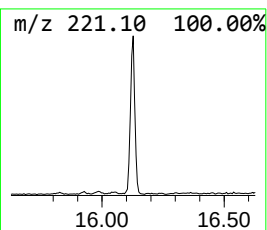
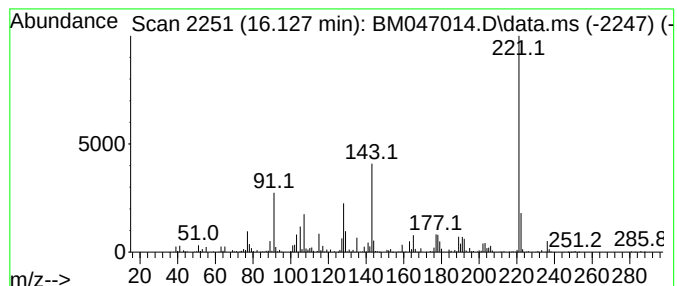
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1H-Indene, 2,3-dihydro-1,1,... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.127	2.73 ng/ul	427328	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	93		
2	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	64		
3	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	43		
4	1H-Indene, 2,3-dihydro-1,1,3-tri...	236	C18H20	003910-35-8	42		
5	2-Oxa-1,3-disilacyclohexane, 1,1...	236	C12H20OSi2	1010308-85-9	38		



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
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Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
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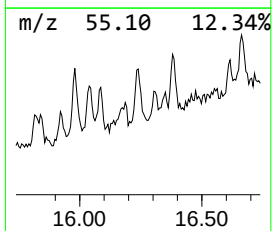
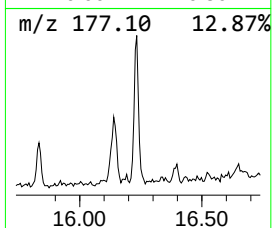
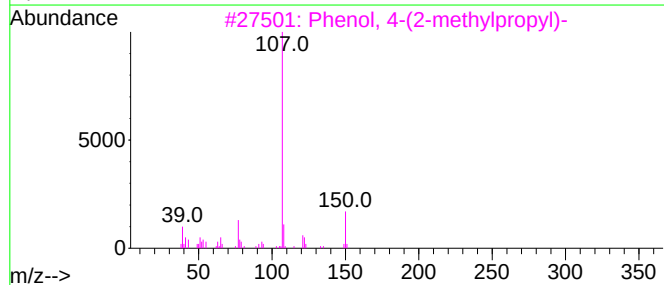
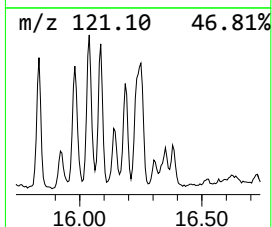
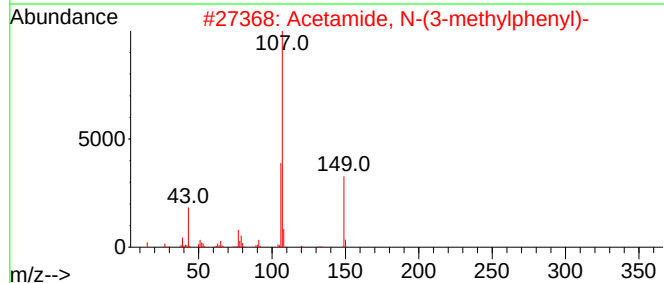
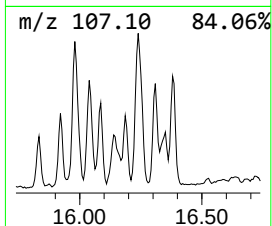
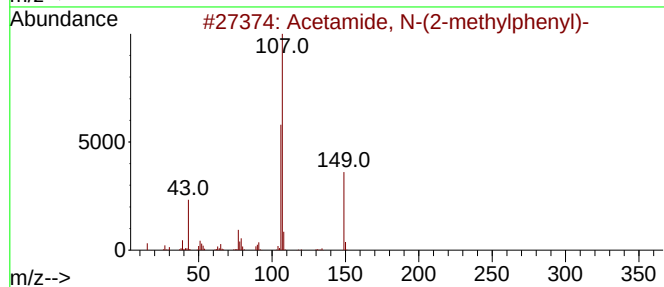
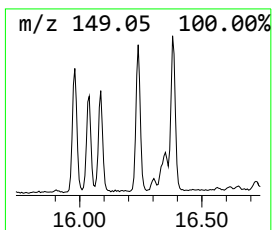
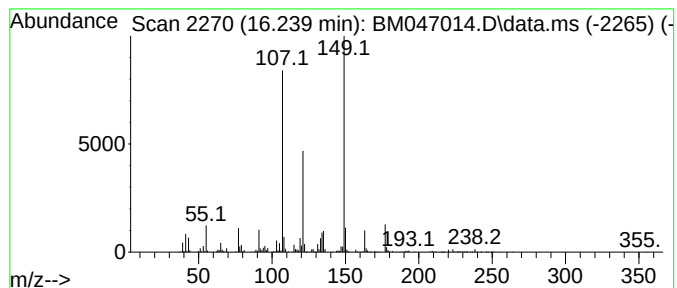
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 unknown-03 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.239	5.62 ng/ul	881113	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Acetamide, N-(2-methylphenyl)-	149	C9H11NO	000120-66-1	46
2			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	46
3			Phenol, 4-(2-methylpropyl)-	150	C10H14O	004167-74-2	42
4			3,4-Methylenedioxypropiofenone	178	C10H10O3	028281-49-4	35
5			Phthalic acid, 5-methylhex-2-yl ...	292	C17H24O4	1000371-09-3	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
Operator : RC/JU
Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
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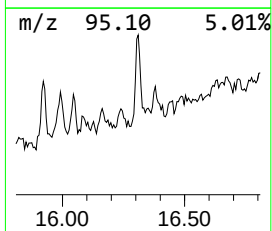
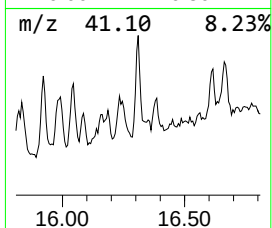
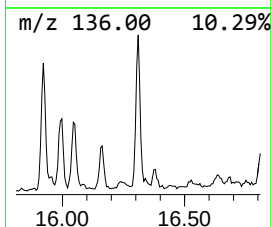
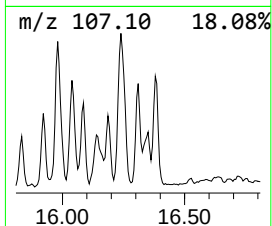
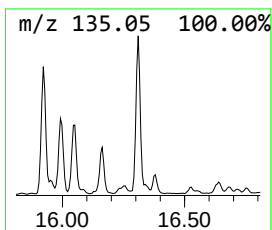
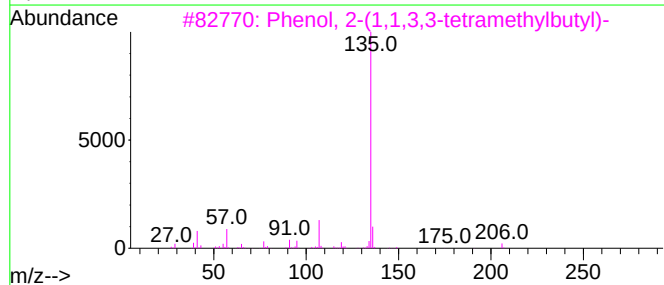
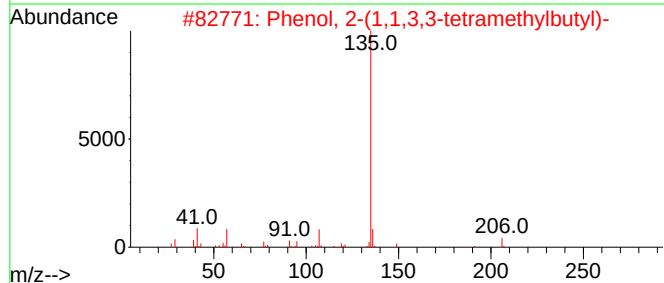
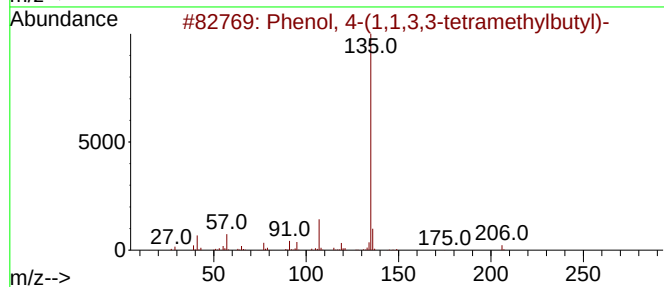
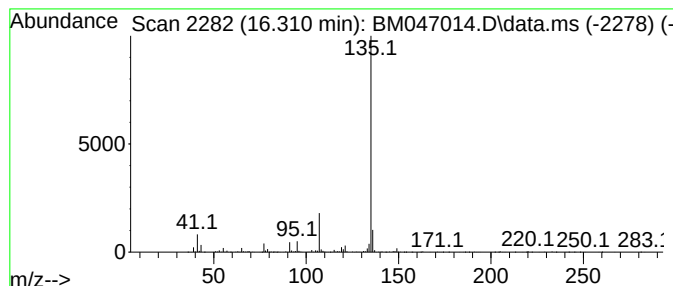
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Phenol, 4-(1,1,3,3-tetramet... Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.310	5.12 ng/ul	802400	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
2			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
3			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	78
4			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78
5			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
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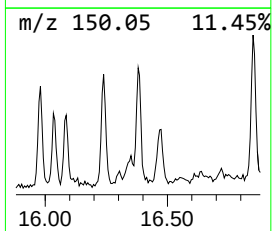
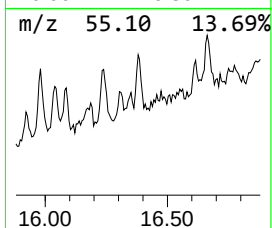
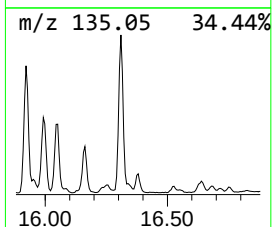
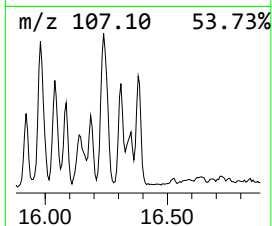
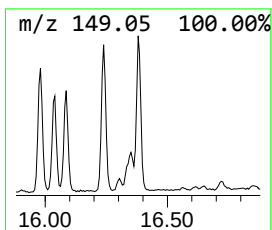
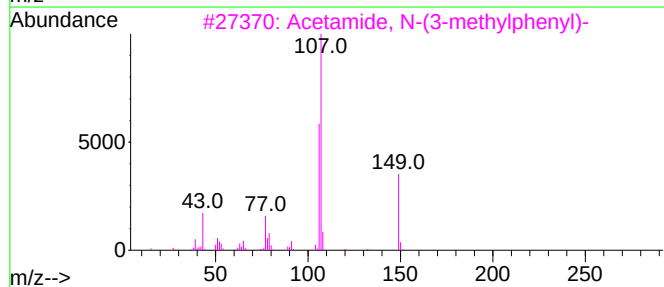
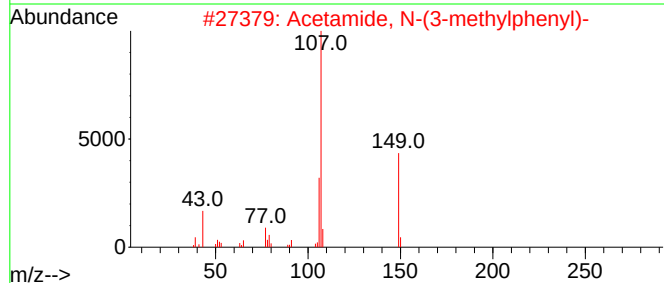
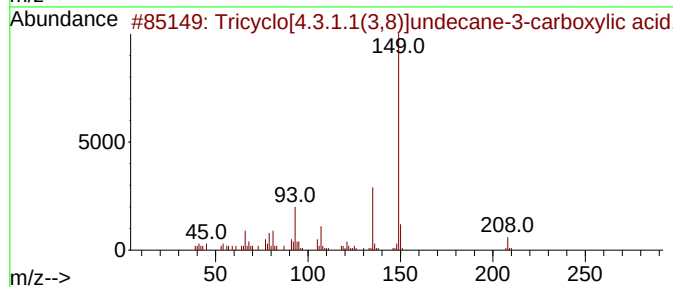
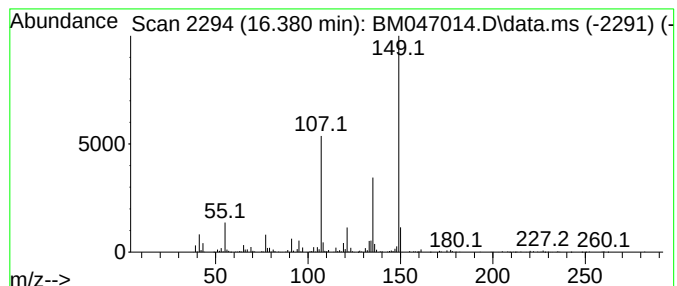
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Tricyclo[4.3.1.1(3,8)]undec... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.380	3.37 ng/ul	528820	Phenanthrene-d10	16.810

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tricyclo[4.3.1.1(3,8)]undecane-3...	208	C13H20O2	031061-61-7	59
2			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	58
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	53
4			2-tert-Butyl-6-methylphenol, n-p...	234	C16H26O	1000395-18-9	47
5			7-Methylthieno[3,2-b]pyridine	149	C8H7NS	013362-83-9	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM080524\
Data File : BM047014.D
Acq On : 06 Aug 2024 06:52
Operator : RC/JU
Sample : P3382-01
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
E20R1

Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-BM080224.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
Benzyl alcohol	7.640	11.4	ng/ul	904245	1	7.393	1587060	20.0
1-Phenoxypropan...	10.910	2.3	ng/ul	241303	2	10.151	2091600	20.0
unknown-01	15.833	3.2	ng/ul	504577	4	16.810	3135770	20.0
Phenol, 2-(1,1,...	15.921	5.1	ng/ul	801903	4	16.810	3135770	20.0
unknown-02	15.980	5.2	ng/ul	813655	4	16.810	3135770	20.0
4-(7-Methylocty...	16.045	4.6	ng/ul	720976	4	16.810	3135770	20.0
1H-Indene, 2,3-...	16.127	2.7	ng/ul	427328	4	16.810	3135770	20.0
unknown-03	16.239	5.6	ng/ul	881113	4	16.810	3135770	20.0
Phenol, 4-(1,1,...	16.310	5.1	ng/ul	802400	4	16.810	3135770	20.0
Tricyclo[4.3.1....	16.380	3.4	ng/ul	528820	4	16.810	3135770	20.0