

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
 Data File : BP018210.D
 Acq On : 16 Nov 2023 13:00
 Operator : MA/JU
 Sample : 05338-15
 Misc :
 ALS Vial : 38 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 GCDM0

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-BP111023.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP018210.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.164	9	13	22	rVB	16440	23480	0.98%	0.099%
2	3.446	59	61	64	rBV4	2368	2628	0.11%	0.011%
3	3.605	85	88	106	rBV	58283	114904	4.79%	0.485%
4	3.770	111	116	117	rBV5	5376	6796	0.28%	0.029%
5	3.899	135	138	146	rVB4	5073	9350	0.39%	0.039%
6	4.417	221	226	233	rVB2	23957	36804	1.53%	0.155%
7	5.034	325	331	337	rBV4	11865	21255	0.89%	0.090%
8	5.623	428	431	436	rBV7	1271	2443	0.10%	0.010%
9	5.846	463	469	475	rBV9	3387	8069	0.34%	0.034%
10	6.240	533	536	540	rVB5	1815	3151	0.13%	0.013%
11	6.422	560	567	572	rBV10	2052	5289	0.22%	0.022%
12	6.964	650	659	670	rBV2	44711	105269	4.39%	0.444%
13	7.134	682	688	700	rBV	203574	327444	13.65%	1.381%
14	7.240	704	706	711	rVV6	1426	2623	0.11%	0.011%
15	7.305	711	717	735	rVV	219401	403478	16.81%	1.702%
16	7.422	735	737	740	rVB4	3087	2668	0.11%	0.011%
17	7.781	792	798	812	rBV	282764	479736	19.99%	2.024%
18	8.516	917	923	947	rBV	140718	308909	12.87%	1.303%
19	8.669	947	949	954	rVB6	2070	2623	0.11%	0.011%
20	8.993	997	1004	1017	rBV	250399	483286	20.14%	2.039%
21	9.158	1030	1032	1038	rVB7	3636	4249	0.18%	0.018%
22	9.316	1054	1059	1060	rBV4	2021	2550	0.11%	0.011%
23	9.705	1118	1125	1141	rBV	234456	459375	19.14%	1.938%
24	9.840	1147	1148	1155	rVB7	2774	3593	0.15%	0.015%
25	10.234	1208	1215	1236	rBV	252954	599854	25.00%	2.531%
26	10.481	1255	1257	1261	rBV5	1854	2753	0.11%	0.012%
27	10.599	1269	1277	1288	rBV	418777	689734	28.74%	2.910%
28	10.793	1303	1310	1332	rBV	117743	333776	13.91%	1.408%
29	10.987	1340	1343	1354	rVB	7770	18222	0.76%	0.077%
30	11.146	1363	1370	1371	rBV6	5880	9858	0.41%	0.042%
31	11.322	1396	1400	1405	rBV8	3544	4500	0.19%	0.019%
32	11.410	1405	1415	1423	rBV6	16148	49930	2.08%	0.211%
33	11.481	1423	1427	1430	rVV5	4544	6691	0.28%	0.028%
34	11.528	1432	1435	1441	rVB6	5753	9997	0.42%	0.042%
35	11.628	1451	1452	1461	rVB9	2536	4178	0.17%	0.018%
36	12.234	1549	1555	1561	rBV9	4030	9642	0.40%	0.041%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
 Title : SVOA CALIBRATION

37	13.299	1730	1736	1749	rVB2	61861	105387	4.39%	0.445%
38	13.675	1793	1800	1805	rBV	249182	408436	17.02%	1.723%
39	13.740	1805	1811	1812	rBV2	27591	49518	2.06%	0.209%
40	13.887	1831	1836	1854	rVB	782341	1133230	47.22%	4.781%
41	14.163	1876	1883	1892	rBV	743984	1081183	45.05%	4.561%
42	14.463	1926	1934	1947	rVB	610416	887979	37.00%	3.746%
43	14.557	1947	1950	1955	rBV6	1404	2586	0.11%	0.011%
44	14.793	1985	1990	2001	rBV6	13437	49884	2.08%	0.210%
45	15.069	2034	2037	2043	rVV8	3444	5237	0.22%	0.022%
46	15.134	2047	2048	2055	rVB7	2039	3568	0.15%	0.015%
47	15.204	2055	2060	2064	rBV6	4010	5736	0.24%	0.024%
48	15.328	2077	2081	2085	rVB7	1919	3099	0.13%	0.013%
49	15.469	2098	2105	2122	rBV	958935	1452904	60.54%	6.129%
50	15.628	2127	2132	2144	rBV	309880	523190	21.80%	2.207%
51	15.881	2171	2175	2179	rBV	26082	36101	1.50%	0.152%
52	15.922	2179	2182	2187	rVV	27073	37745	1.57%	0.159%
53	15.963	2187	2189	2194	rVV2	7734	9967	0.42%	0.042%
54	16.010	2194	2197	2200	rVB4	2864	3949	0.16%	0.017%
55	16.093	2206	2211	2213	rBV6	2955	4864	0.20%	0.021%
56	16.181	2222	2226	2235	rVB	45195	66570	2.77%	0.281%
57	16.275	2235	2242	2247	rBV	150841	218883	9.12%	0.923%
58	16.334	2247	2252	2258	rVV3	190615	389629	16.24%	1.644%
59	16.398	2258	2263	2267	rVV2	157961	278295	11.60%	1.174%
60	16.440	2267	2270	2275	rVV	95156	134288	5.60%	0.567%
61	16.498	2275	2280	2282	rVV2	53503	87620	3.65%	0.370%
62	16.545	2286	2288	2292	rVV	59525	68717	2.86%	0.290%
63	16.598	2292	2297	2304	rVV3	129122	255349	10.64%	1.077%
64	16.669	2304	2309	2314	rVV	155138	243106	10.13%	1.026%
65	16.716	2314	2317	2319	rVV2	49677	70465	2.94%	0.297%
66	16.745	2319	2322	2329	rVB	87847	114514	4.77%	0.483%
67	16.881	2342	2345	2348	rVB6	4881	6047	0.25%	0.026%
68	16.916	2349	2351	2356	rVB6	3473	4223	0.18%	0.018%
69	16.963	2356	2359	2360	rBV3	2805	3151	0.13%	0.013%
70	17.004	2362	2366	2369	rVV2	6334	10106	0.42%	0.043%
71	17.045	2370	2373	2376	rVV5	3907	5075	0.21%	0.021%
72	17.081	2376	2379	2383	rVV6	6767	9414	0.39%	0.040%
73	17.110	2383	2384	2390	rVB4	4044	5282	0.22%	0.022%
74	17.245	2395	2407	2417	rBV	853528	1192924	49.71%	5.033%
75	17.345	2418	2424	2442	rVB	1244607	1803480	75.15%	7.608%
76	17.522	2452	2454	2459	rVB2	10472	12256	0.51%	0.052%

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77	17.581	2459	2464	2467	rBV2	16807	22311	0.93%	0.094%
78	17.610	2467	2469	2476	rVB6	7931	12565	0.52%	0.053%
79	17.669	2476	2479	2483	rBV2	12921	16395	0.68%	0.069%
80	17.710	2484	2486	2490	rVB4	4913	4837	0.20%	0.020%
81	17.751	2490	2493	2497	rVB6	10128	12267	0.51%	0.052%
82	17.798	2497	2501	2504	rBV3	13894	19046	0.79%	0.080%
83	17.845	2507	2509	2510	rVV2	5063	4219	0.18%	0.018%
84	17.881	2510	2515	2525	rVV2	384215	486060	20.25%	2.051%
85	17.963	2527	2529	2533	rVB4	7119	6884	0.29%	0.029%
86	18.040	2539	2542	2546	rBV6	6316	11517	0.48%	0.049%
87	18.098	2548	2552	2559	rBV2	39035	60467	2.52%	0.255%
88	18.192	2564	2568	2572	rVB5	15657	22251	0.93%	0.094%
89	18.310	2584	2588	2593	rBV	67245	88521	3.69%	0.373%
90	18.669	2646	2649	2655	rBV8	5687	8598	0.36%	0.036%
91	19.004	2703	2706	2713	rBV8	7480	15464	0.64%	0.065%
92	19.175	2731	2735	2739	rBV4	17298	21238	0.89%	0.090%
93	19.234	2742	2745	2746	rVV3	10265	12140	0.51%	0.051%
94	19.257	2746	2749	2756	rVV2	17026	27881	1.16%	0.118%
95	19.345	2760	2764	2774	rVV7	20366	34904	1.45%	0.147%
96	19.481	2784	2787	2792	rVB7	8645	10534	0.44%	0.044%
97	19.604	2802	2808	2817	rBV	1801907	2324192	96.85%	9.805%
98	21.381	3105	3110	3119	rBV	990533	1334151	55.60%	5.628%
99	23.686	3495	3502	3517	rVB2	1196316	2399723	100.00%	10.124%
100	23.851	3523	3530	3543	rVB	681716	1404751	58.54%	5.926%

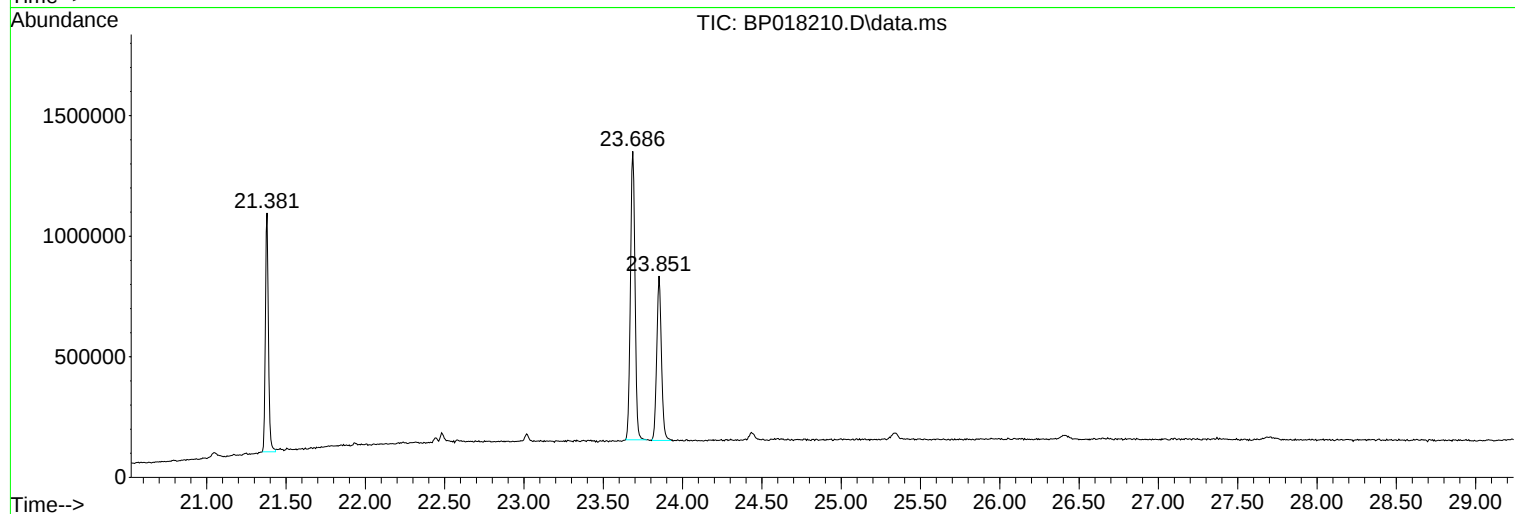
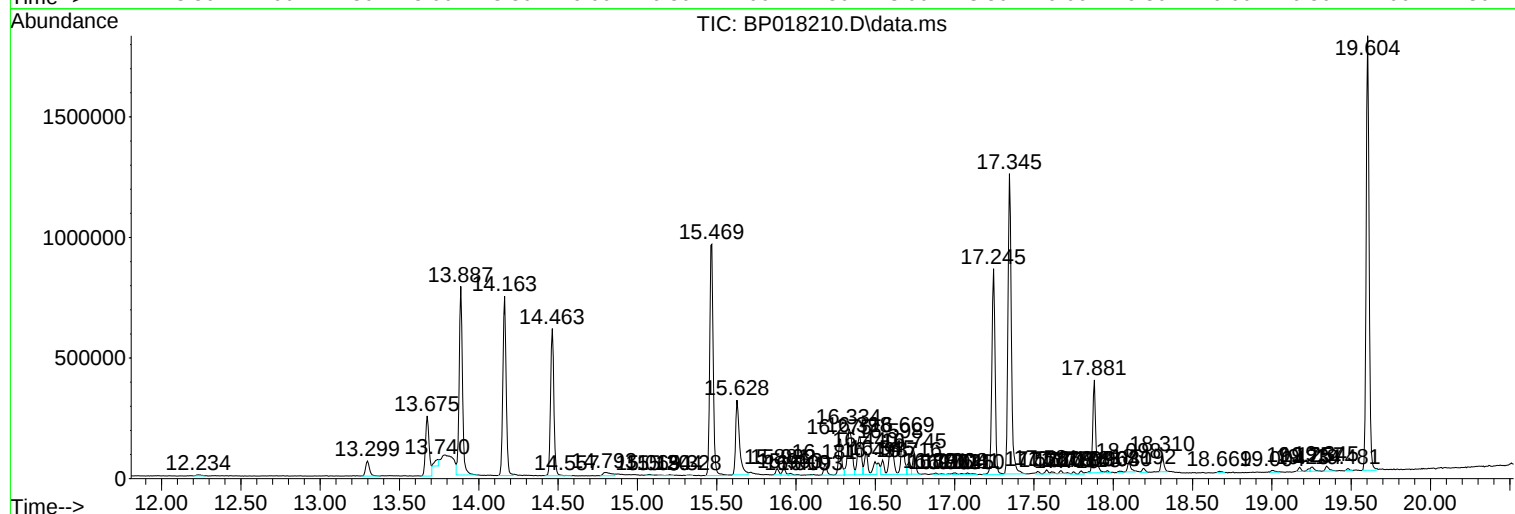
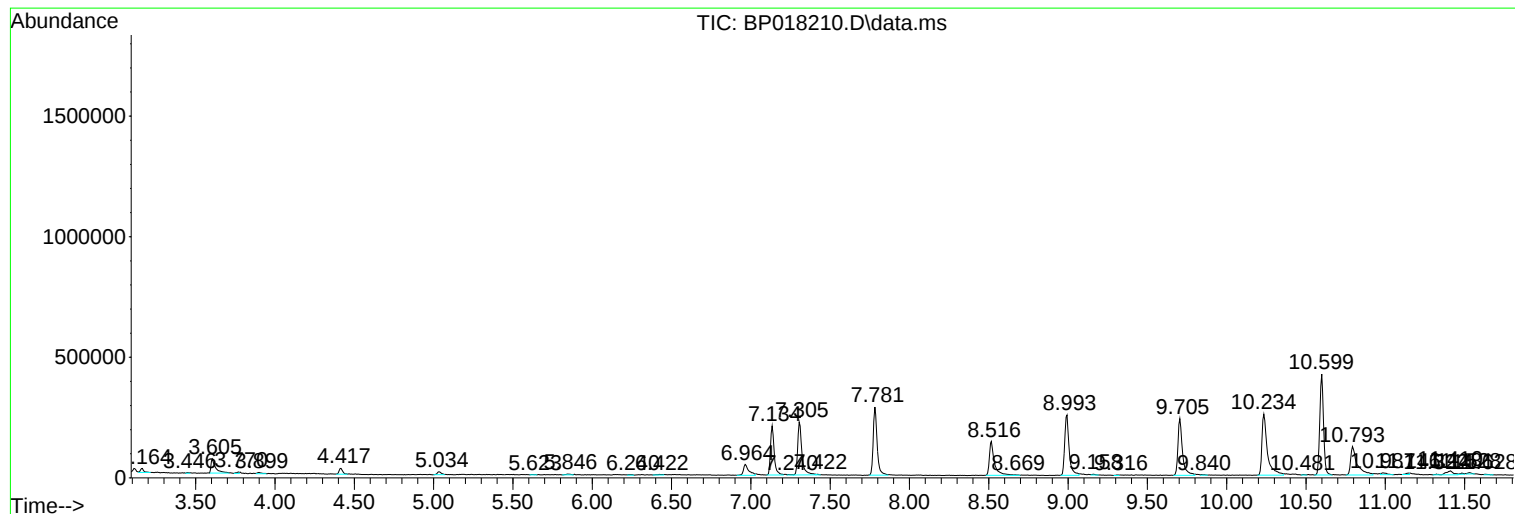
Sum of corrected areas: 23703980

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BNA_P
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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



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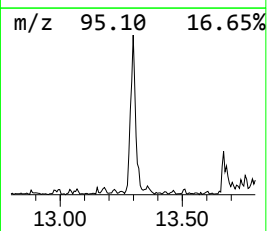
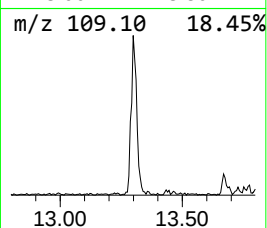
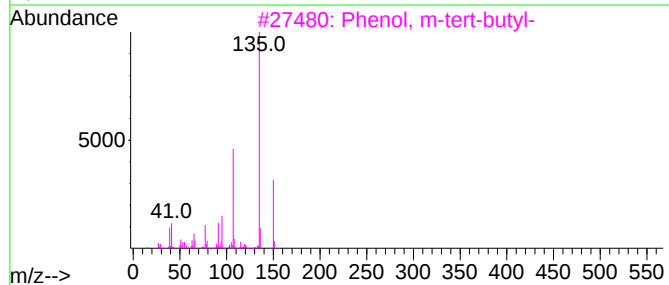
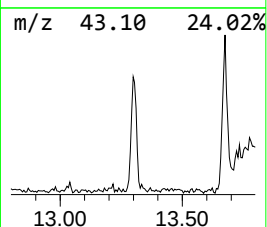
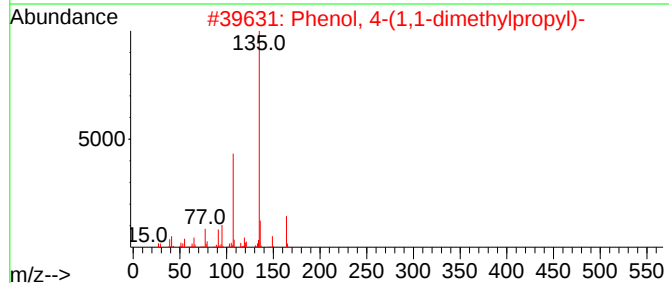
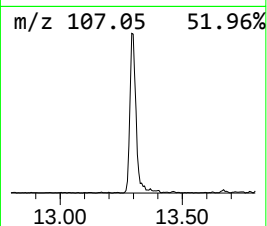
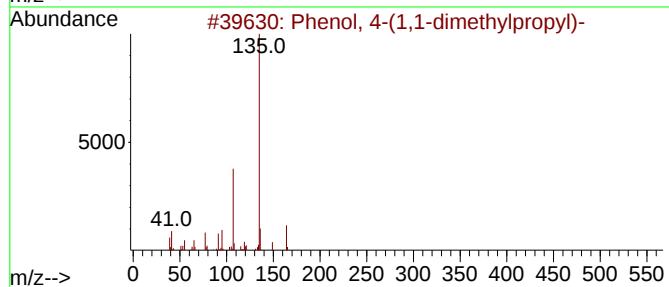
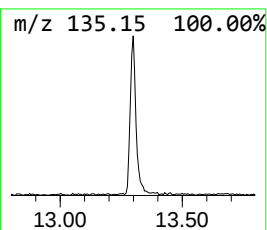
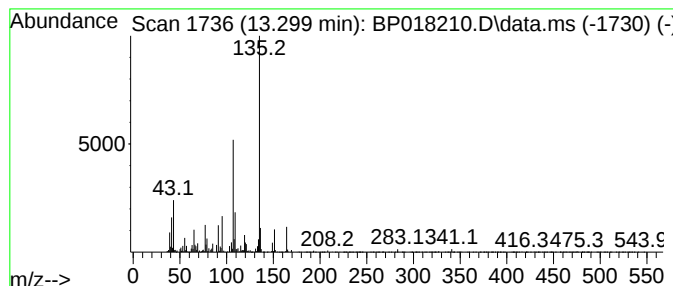
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Phenol, 4-(1,1-dimethylprop... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.299	2.37 ng/ul	105387	Acenaphthene-d10	14.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	93
2			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	91
3			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
4			Phenol, m-tert-butyl-	150	C10H14O	000585-34-2	64
5			Propane, 1,3-bis(ethylthio)-	164	C7H16S2	033672-52-5	58



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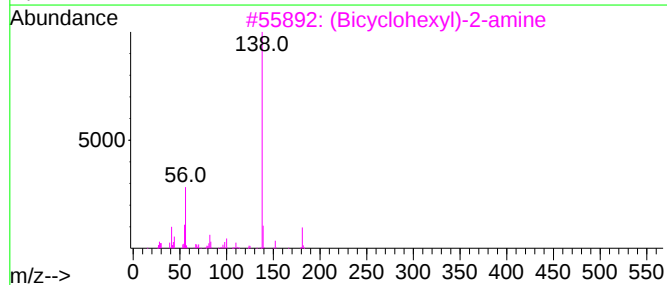
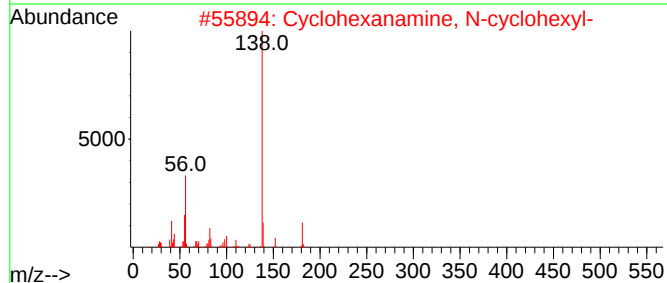
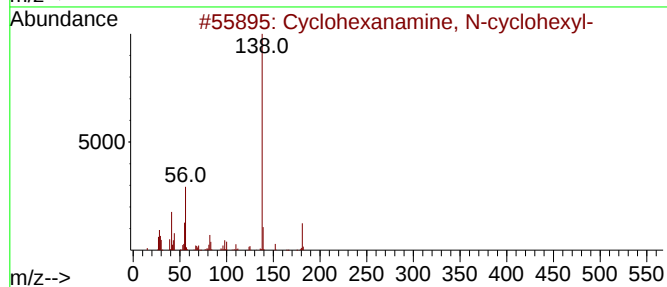
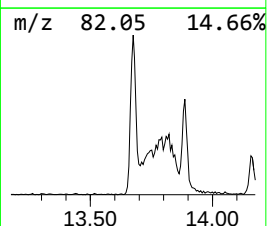
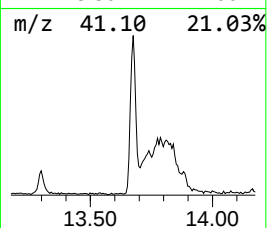
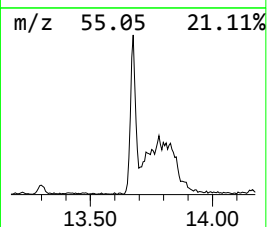
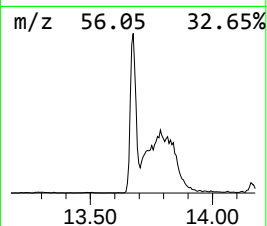
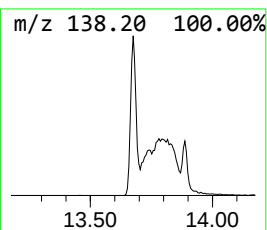
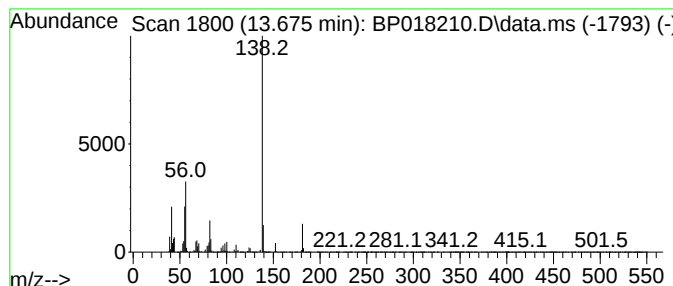
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Cyclohexanamine, N-cyclohexyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.675	9.20 ng/ul	408436	Acenaphthene-d10	14.463

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	97
2			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	96
3			(Bicyclohexyl)-2-amine	181	C12H23N	006283-14-3	95
4			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	94
5			Cyclohexanamine, N-cyclohexyl-	181	C12H23N	000101-83-7	91



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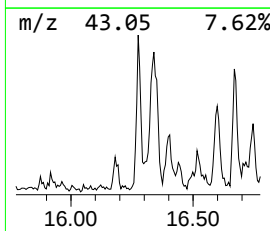
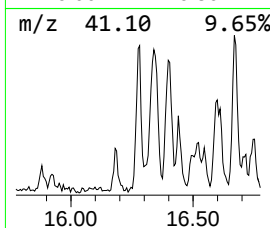
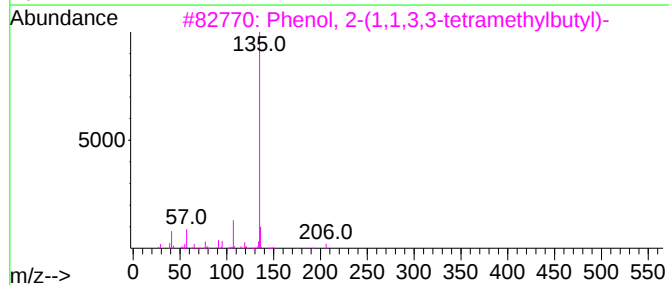
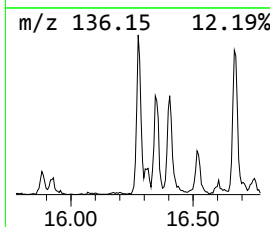
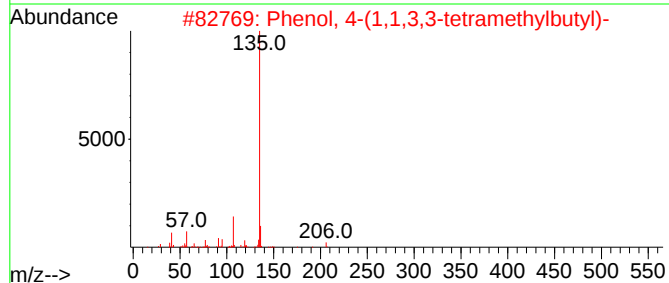
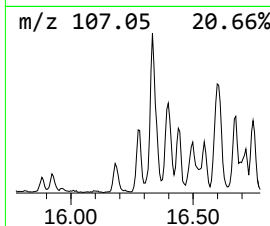
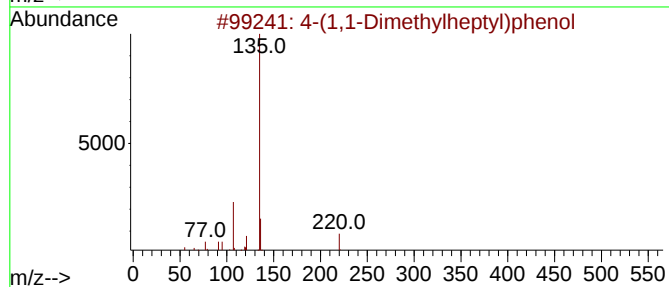
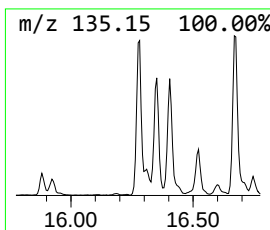
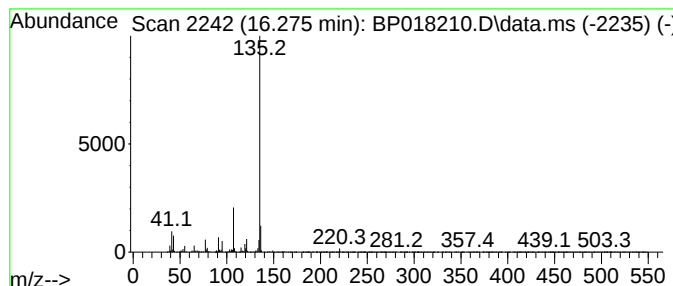
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Quant Title : SVOA CALIBRATION

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

Peak Number 3 4-(1,1-Dimethylheptyl)phenol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.275	3.67 ng/ul	218883	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(1,1-Dimethylheptyl)phenol	220	C15H24O	1000384-80-3	91
2			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	86
3			Phenol, 2-(1,1,3,3-tetramethylbu...	206	C14H22O	003884-95-5	80
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
5			4-tert-Butylphenyl acetate	192	C12H16O2	003056-64-2	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDM0

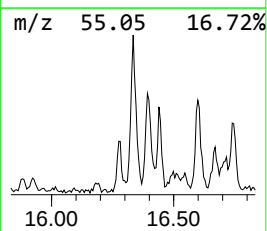
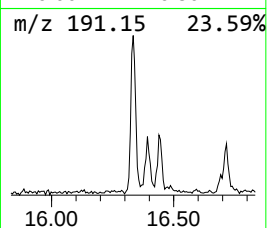
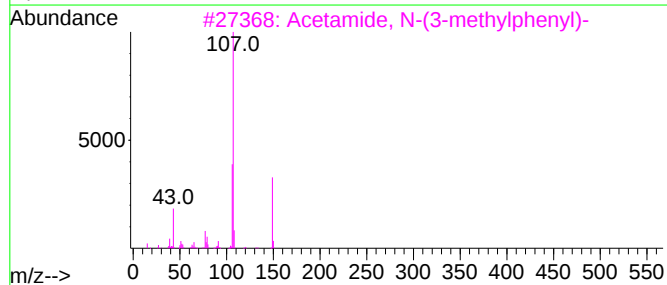
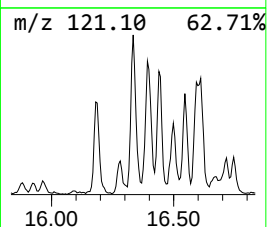
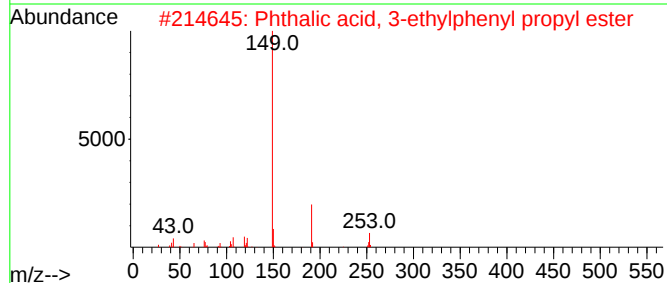
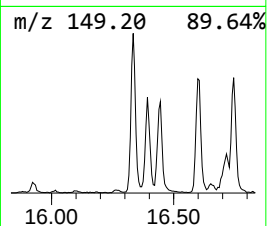
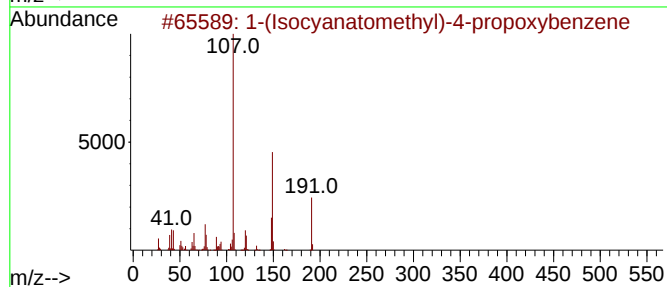
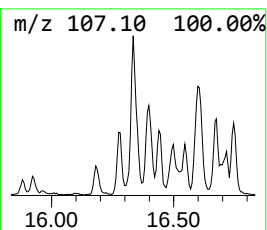
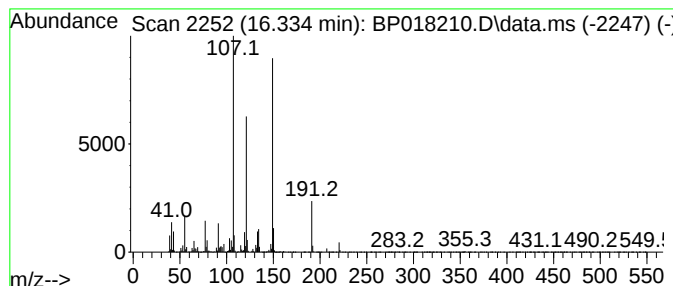
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 1-(Isocyanatomethyl)-4-prop... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.334	6.53 ng/ul	389629	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-(Isocyanatomethyl)-4-propoxybe...	191	C11H13NO2	936095-07-7	72
2			Phthalic acid, 3-ethylphenyl pro...	312	C19H20O4	1000357-07-6	43
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
4			Phthalic acid, 3-fluorophenyl pr...	302	C17H15FO4	1000356-49-8	38
5			Ethanol, 2-(3,3-dimethylbicyclo[...	166	C11H18O	002226-05-3	38



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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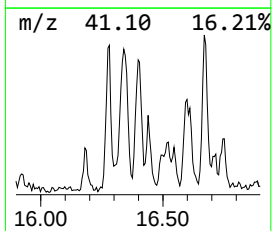
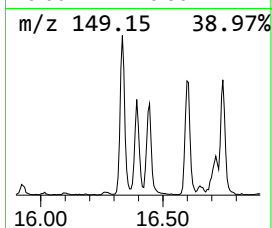
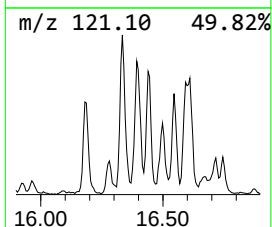
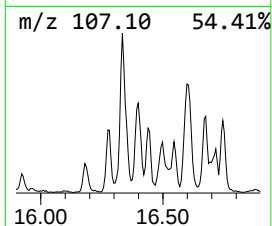
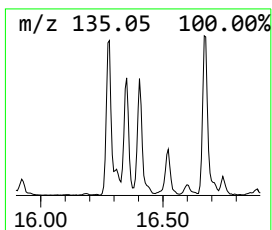
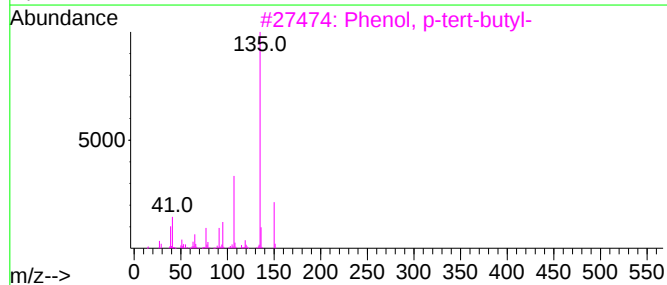
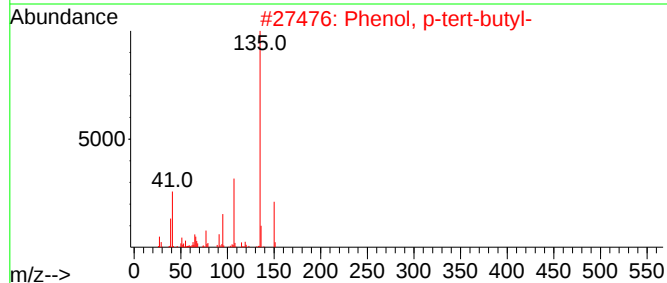
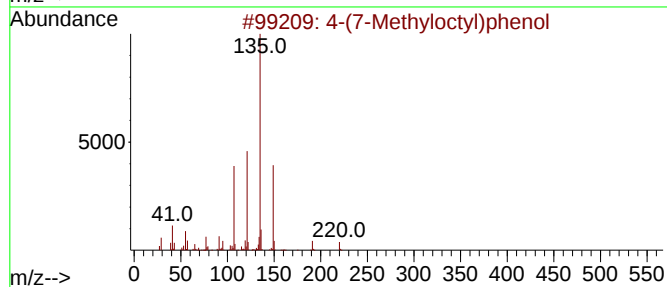
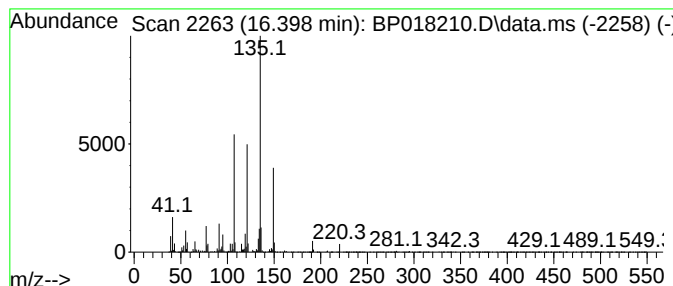
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 4-(7-Methyloctyl)phenol Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.398	4.67 ng/ul	278295	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(7-Methyloctyl)phenol	220	C15H24O	024518-48-7	98
2			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
3			Phenol, p-tert-butyl-	150	C10H14O	000098-54-4	64
4			Hexestrol, 0-methoxycarbonyl-	328	C20H24O4	1000487-66-3	53
5			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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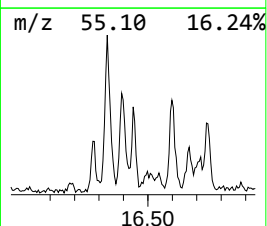
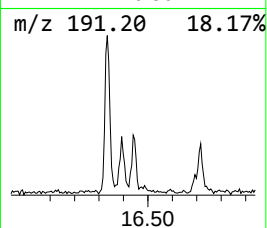
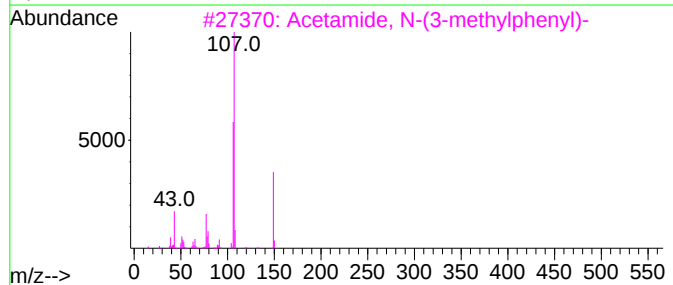
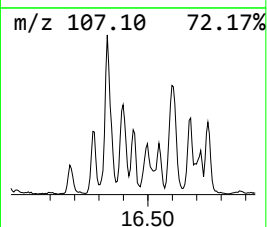
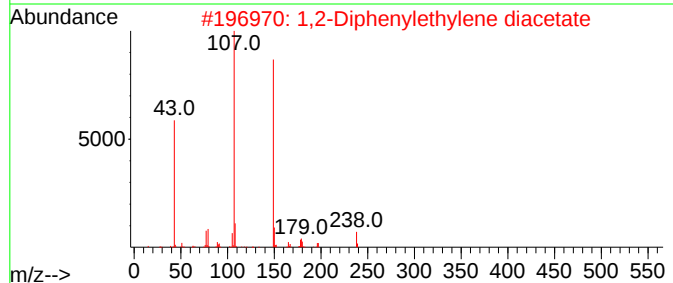
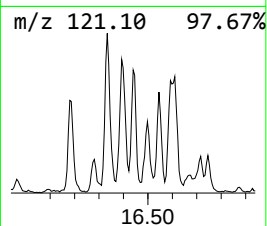
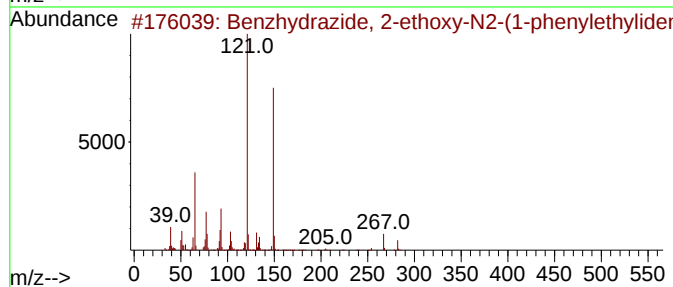
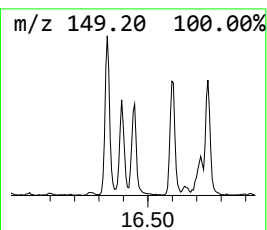
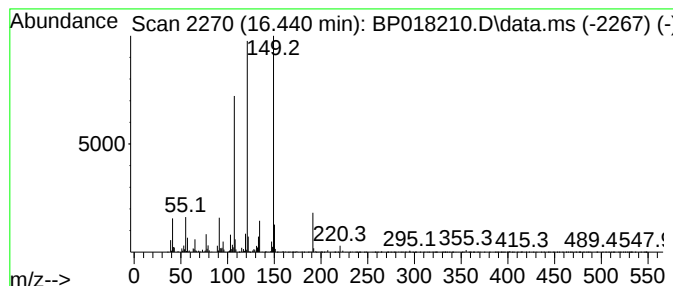
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzhydrazide, 2-ethoxy-N2-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.440	2.25 ng/ul	134288	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzhydrazide, 2-ethoxy-N2-(1-ph...	282	C17H18N2O2	331751-10-1	59
2			1,2-Diphenylethylene diacetate	298	C18H18O4	003682-07-3	43
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	43
4			Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	35
5			Phenol, 4-(1-methylpropyl)-	150	C10H14O	000099-71-8	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
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Instrument :
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ClientSampleId :
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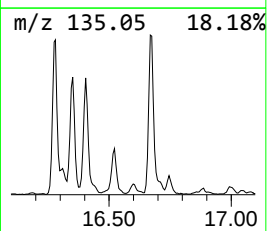
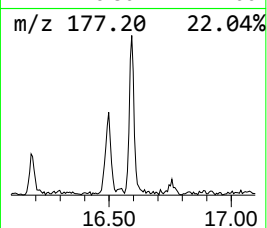
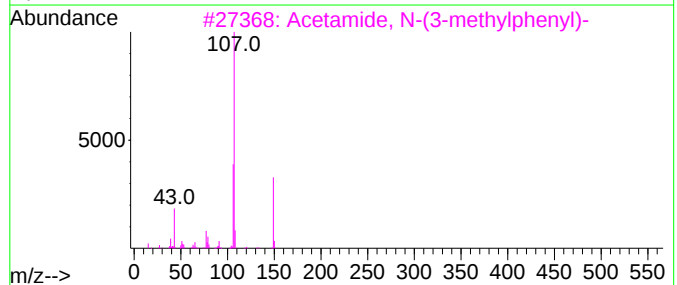
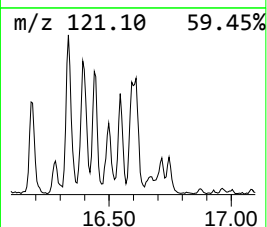
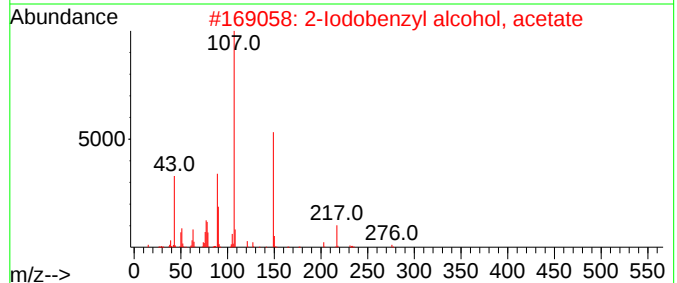
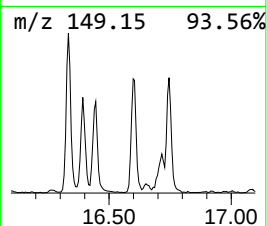
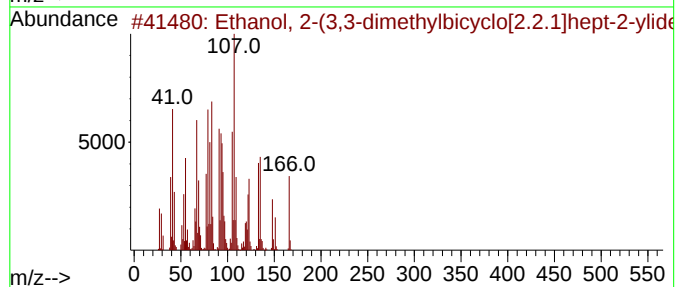
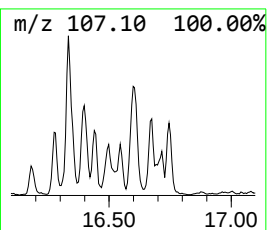
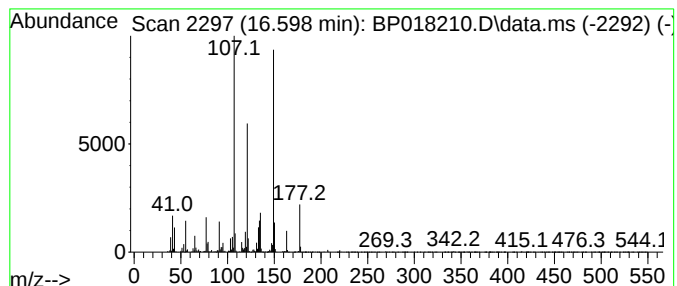
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Ethanol, 2-(3,3-dimethylbic... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.598	4.28 ng/ul	255349	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethanol, 2-(3,3-dimethylbicyclo[...]	166	C11H18O	002226-05-3	80
2			2-Iodobenzyl alcohol, acetate	276	C9H9IO2	1000364-47-1	50
3			Acetamide, N-(3-methylphenyl)-	149	C9H11NO	000537-92-8	38
4			Methyl (1s,3s,5R,7S)-3-methylada...	208	C13H20O2	1010504-39-1	35
5			Benzene, 1-ethoxy-4-ethyl-	150	C10H14O	001585-06-4	30



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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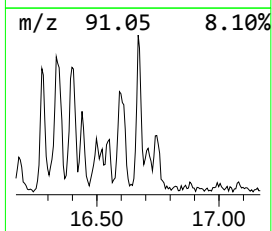
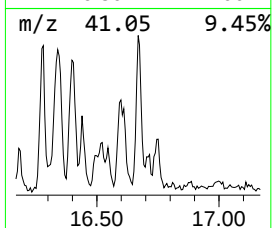
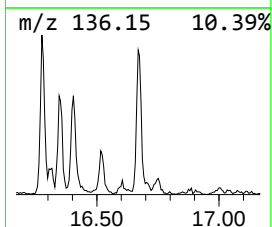
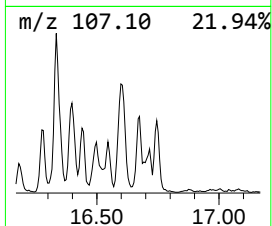
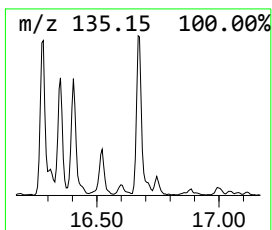
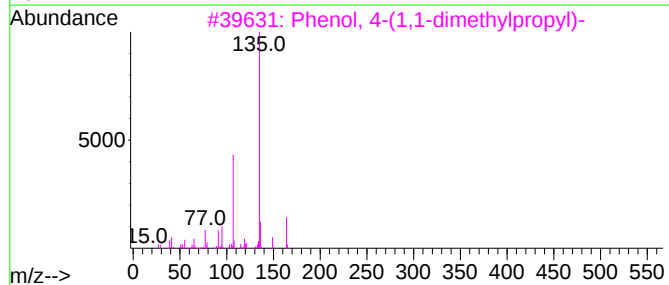
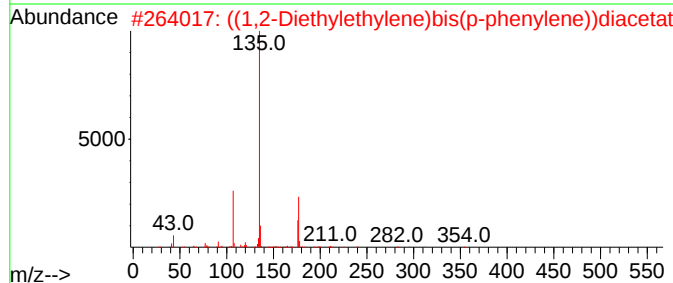
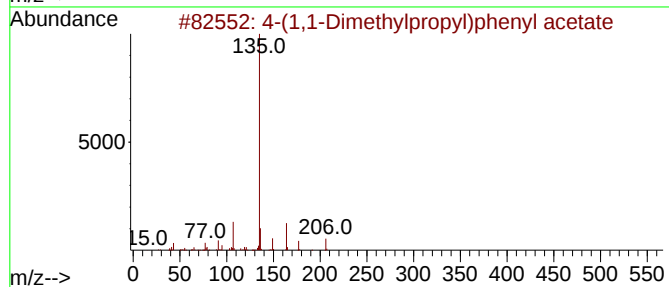
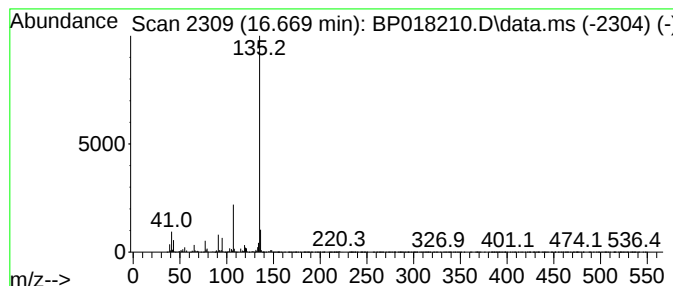
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 4-(1,1-Dimethylpropyl)pheny... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.669	4.08 ng/ul	243106	Phenanthrene-d10	17.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			4-(1,1-Dimethylpropyl)phenyl ace...	206	C13H18O2	1000373-45-3	83
2			((1,2-Diethylethylene)bis(p-phen...	354	C22H26O4	004547-76-6	78
3			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
4			Phenol, 4-(1,1-dimethylpropyl)-	164	C11H16O	000080-46-6	78
5			Phenol, 4-(1,1,3,3-tetramethylbu...	206	C14H22O	000140-66-9	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
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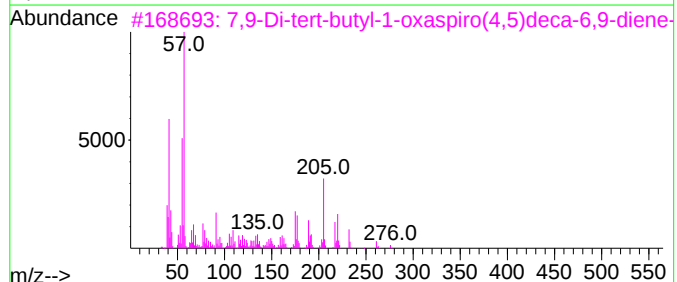
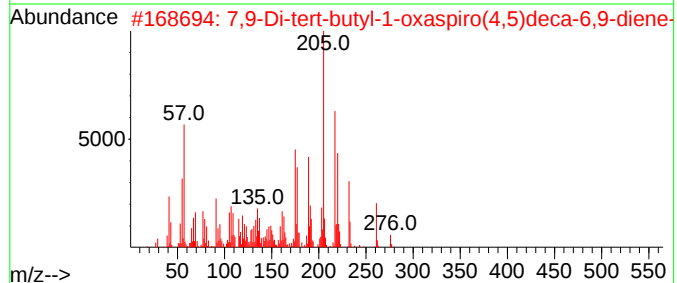
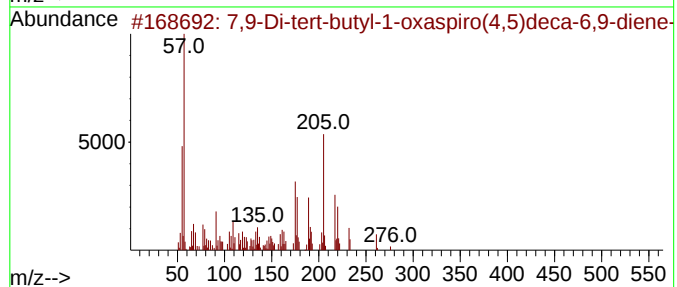
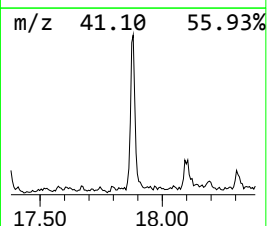
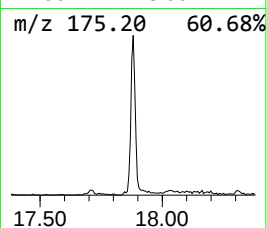
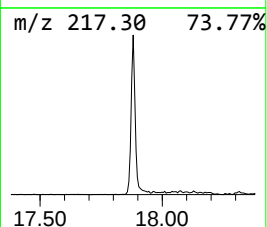
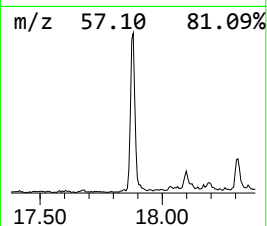
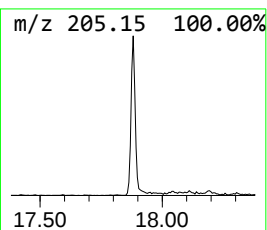
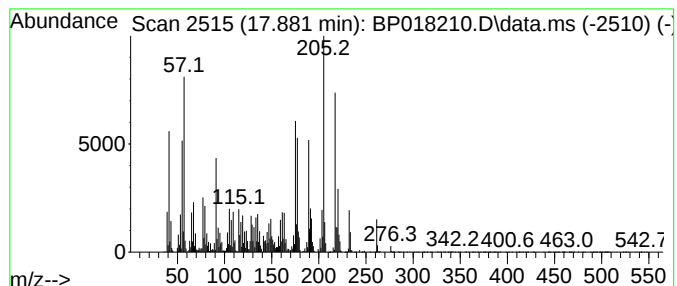
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ClientSampleId :
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 7,9-Di-tert-butyl-1-oxaspiro... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
17.881	8.15 ng/ul	486060	Phenanthrene-d10	17.245	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	99
2	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	97
3	7,9-Di-tert-butyl-1-oxaspiro(4,5...	276	C17H24O3	082304-66-3	50
4	Ethanone, 1-(5,6,7,8-tetrahydro-...	220	C14H20O2	071596-88-8	42
5	1-(3-Amino-4,6-dimethylthieno[2,...	220	C11H12N2OS	052505-42-7	41



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP111523\
Data File : BP018210.D
Acq On : 16 Nov 2023 13:00
Operator : MA/JU
Sample : 05338-15
Misc :
ALS Vial : 38 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
GCDM0

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP111023.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
Phenol, 4-(1,1-...	13.299	2.4	ng/ul	105387	3	14.463	887979	20.0
Cyclohexanamine...	13.675	9.2	ng/ul	408436	3	14.463	887979	20.0
4-(1,1-Dimethyl...	16.275	3.7	ng/ul	218883	4	17.245	1192920	20.0
1-(Isocyanatome...	16.334	6.5	ng/ul	389629	4	17.245	1192920	20.0
4-(7-Methylocty...	16.398	4.7	ng/ul	278295	4	17.245	1192920	20.0
Benzhydrazide, ...	16.440	2.3	ng/ul	134288	4	17.245	1192920	20.0
Ethanol, 2-(3,3...	16.598	4.3	ng/ul	255349	4	17.245	1192920	20.0
4-(1,1-Dimethyl...	16.669	4.1	ng/ul	243106	4	17.245	1192920	20.0
7,9-Di-tert-but...	17.881	8.2	ng/ul	486060	4	17.245	1192920	20.0