

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
 Data File : BP010547.D
 Acq On : 25 May 2022 21:31
 Operator : CG/JU
 Sample : N2950-10DL 5X
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 DBTD0DL

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-BP051322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010547.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.781	114	118	129	rBV2	24200	59568	2.52%	0.215%
2	5.393	388	392	400	rVV5	7135	12876	0.54%	0.046%
3	5.528	409	415	422	rBV3	15295	32388	1.37%	0.117%
4	5.899	473	478	484	rVB2	11781	18682	0.79%	0.067%
5	6.975	651	661	666	rVV	16633	29045	1.23%	0.105%
6	7.052	668	674	680	rVV3	19658	46019	1.95%	0.166%
7	7.111	680	684	689	rVB	14556	24252	1.03%	0.087%
8	7.216	696	702	709	rBV	93616	164879	6.98%	0.595%
9	7.416	728	736	745	rBV	84856	148101	6.27%	0.534%
10	7.534	750	756	765	rVV	37115	64747	2.74%	0.234%
11	7.881	807	815	827	rBV	477021	799913	33.85%	2.886%
12	8.005	830	836	842	rVB2	18009	33908	1.43%	0.122%
13	8.258	871	879	885	rVV2	50703	86994	3.68%	0.314%
14	8.422	902	907	916	rVB4	12724	25366	1.07%	0.092%
15	8.528	921	925	930	rVV4	8441	15844	0.67%	0.057%
16	8.599	930	937	944	rVV	60755	118689	5.02%	0.428%
17	8.687	947	952	956	rVV7	9947	25257	1.07%	0.091%
18	8.734	956	960	966	rVV6	11919	25218	1.07%	0.091%
19	8.993	999	1004	1007	rBV3	10962	18059	0.76%	0.065%
20	9.046	1007	1013	1022	rVV2	113816	231462	9.79%	0.835%
21	9.240	1041	1046	1052	rVB2	12515	18468	0.78%	0.067%
22	9.346	1055	1064	1066	rBV9	9012	19496	0.82%	0.070%
23	9.575	1099	1103	1108	rVB	9475	14119	0.60%	0.051%
24	9.687	1116	1122	1127	rBV3	11852	21790	0.92%	0.079%
25	9.769	1127	1136	1143	rVV	88889	177355	7.50%	0.640%
26	9.828	1143	1146	1151	rVV4	14053	23914	1.01%	0.086%
27	9.881	1151	1155	1159	rVV7	6755	12880	0.55%	0.046%
28	9.934	1159	1164	1172	rVB5	15346	30624	1.30%	0.110%
29	10.087	1179	1190	1199	rBV2	45830	96259	4.07%	0.347%
30	10.181	1199	1206	1219	rVV4	18884	59169	2.50%	0.213%
31	10.310	1221	1228	1238	rVV3	118418	237315	10.04%	0.856%
32	10.599	1271	1277	1282	rBV8	8555	19537	0.83%	0.070%
33	10.681	1282	1291	1300	rBV	673485	1141645	48.31%	4.119%
34	10.852	1309	1320	1333	rVB2	198063	508644	21.52%	1.835%
35	11.046	1346	1353	1367	rBV	208230	430447	18.21%	1.553%
36	11.299	1391	1396	1401	rVV6	18861	37343	1.58%	0.135%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
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37	11.799	1475	1481	1486	rBV6	10129	18620	0.79%	0.067%
38	12.075	1522	1528	1535	rBV10	15918	34635	1.47%	0.125%
39	12.234	1546	1555	1561	rBV	29296	57903	2.45%	0.209%
40	12.340	1565	1573	1581	rBV	219140	360314	15.25%	1.300%
41	12.457	1589	1593	1599	rVB3	12557	21731	0.92%	0.078%
42	12.557	1599	1610	1623	rBV	561460	993729	42.05%	3.585%
43	12.987	1677	1683	1689	rBV4	20652	38466	1.63%	0.139%
44	13.051	1691	1694	1701	rVB8	13819	25730	1.09%	0.093%
45	13.251	1716	1728	1739	rBV	83311	210408	8.90%	0.759%
46	13.351	1739	1745	1754	rVB	298058	481484	20.37%	1.737%
47	13.498	1765	1770	1774	rBV3	23440	40959	1.73%	0.148%
48	13.546	1775	1778	1779	rBV	17839	16075	0.68%	0.058%
49	13.693	1790	1803	1812	rBV4	68432	200842	8.50%	0.725%
50	13.757	1812	1814	1820	rVV7	8543	13377	0.57%	0.048%
51	13.834	1820	1827	1832	rVV	100994	188872	7.99%	0.681%
52	13.893	1832	1837	1838	rVV	58327	80700	3.41%	0.291%
53	13.922	1838	1842	1852	rVB	348355	508226	21.51%	1.834%
54	14.069	1863	1867	1871	rVV	29178	48752	2.06%	0.176%
55	14.104	1871	1873	1877	rVB4	18534	20924	0.89%	0.075%
56	14.204	1882	1890	1903	rBV	320918	568972	24.08%	2.053%
57	14.510	1933	1942	1948	rBV	1055046	1678915	71.04%	6.057%
58	14.575	1948	1953	1960	rVV	1595033	2363182	100.00%	8.526%
59	14.645	1960	1965	1975	rVB	345604	539221	22.82%	1.945%
60	14.745	1980	1982	1985	rVV4	10867	14177	0.60%	0.051%
61	14.787	1985	1989	1992	rVV6	16654	30179	1.28%	0.109%
62	14.822	1993	1995	2000	rVV3	15427	20428	0.86%	0.074%
63	14.910	2004	2010	2021	rVV2	486808	929871	39.35%	3.355%
64	15.063	2031	2036	2042	rBV	78784	124383	5.26%	0.449%
65	15.145	2042	2050	2055	rVV5	26361	48305	2.04%	0.174%
66	15.440	2087	2100	2106	rBV3	430219	868071	36.73%	3.132%
67	15.504	2106	2111	2116	rVV2	645839	1090631	46.15%	3.935%
68	15.563	2116	2121	2129	rVV	498348	790399	33.45%	2.852%
69	15.634	2129	2133	2139	rVV2	132834	254994	10.79%	0.920%
70	15.687	2139	2142	2145	rVV3	48609	76790	3.25%	0.277%
71	15.728	2145	2149	2156	rVV7	34220	84754	3.59%	0.306%
72	15.787	2156	2159	2161	rVV4	16523	25684	1.09%	0.093%
73	15.816	2161	2164	2167	rVV4	18672	29919	1.27%	0.108%
74	15.857	2167	2171	2176	rVV6	19601	35761	1.51%	0.129%
75	15.934	2180	2184	2191	rVV7	14160	39061	1.65%	0.141%
76	16.004	2191	2196	2209	rVV5	43186	114284	4.84%	0.412%

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 Title : SVOA CALIBRATION

77	16.234	2229	2235	2250	rVB2	129042	247371	10.47%	0.892%
78	16.416	2262	2266	2270	rBV3	11816	20168	0.85%	0.073%
79	16.528	2280	2285	2291	rBV7	18625	43943	1.86%	0.159%
80	16.922	2347	2352	2362	rVB2	73546	122309	5.18%	0.441%
81	17.081	2377	2379	2385	rVB3	13775	19180	0.81%	0.069%
82	17.222	2400	2403	2404	rVV3	14121	12784	0.54%	0.046%
83	17.263	2404	2410	2415	rVV	1424317	2089872	88.43%	7.540%
84	17.304	2415	2417	2421	rVV	108198	153344	6.49%	0.553%
85	17.363	2422	2427	2442	rVB	513076	819017	34.66%	2.955%
86	17.669	2470	2479	2490	rBV	792105	1129370	47.79%	4.075%
87	17.763	2492	2495	2501	rVV6	15744	33464	1.42%	0.121%
88	17.834	2504	2507	2514	rVB2	12834	24047	1.02%	0.087%
89	17.969	2528	2530	2535	rVB4	13140	17943	0.76%	0.065%
90	18.098	2548	2552	2557	rBV4	15226	22391	0.95%	0.081%
91	18.151	2557	2561	2563	rBV4	34481	46876	1.98%	0.169%
92	18.181	2563	2566	2571	rVB	34207	50666	2.14%	0.183%
93	18.469	2610	2615	2621	rVV4	13383	24871	1.05%	0.090%
94	18.557	2626	2630	2637	rVB3	16222	26077	1.10%	0.094%
95	19.275	2749	2752	2755	rVB	15518	17141	0.73%	0.062%
96	19.386	2767	2771	2774	rBV6	11458	19465	0.82%	0.070%
97	19.598	2801	2807	2819	rBV	668759	935156	39.57%	3.374%
98	21.351	3099	3105	3117	rBV	1338073	1945613	82.33%	7.019%
99	23.610	3484	3489	3495	rBV2	233123	467848	19.80%	1.688%
100	23.769	3509	3516	3530	rVB2	684063	1508409	63.83%	5.442%

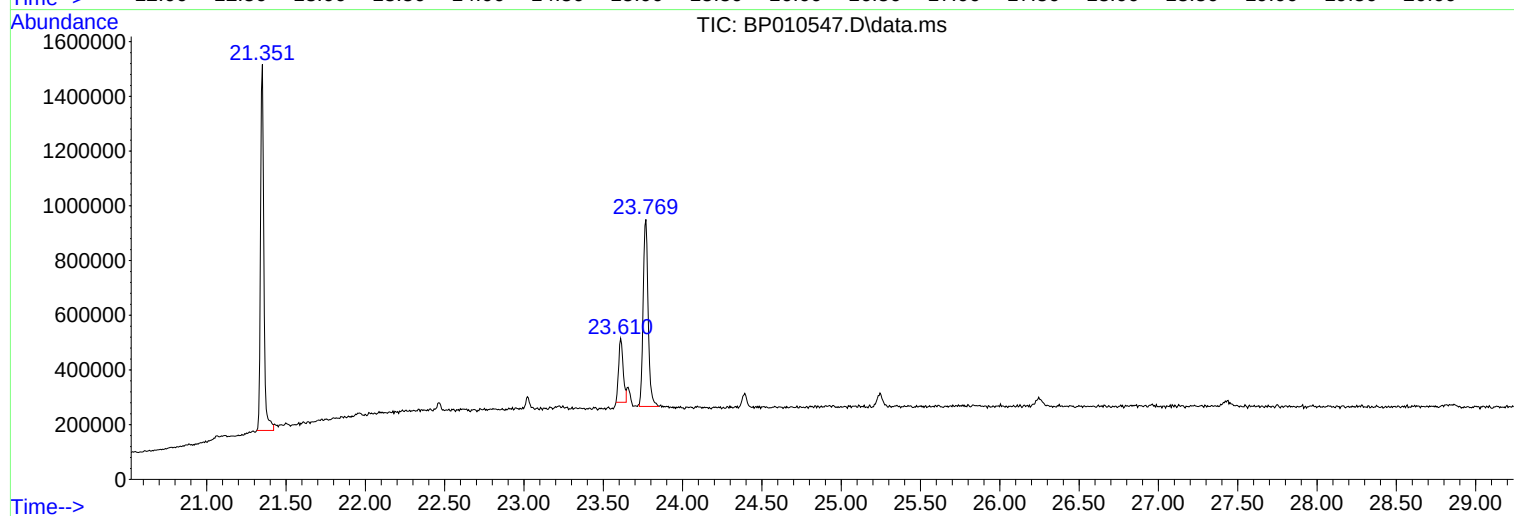
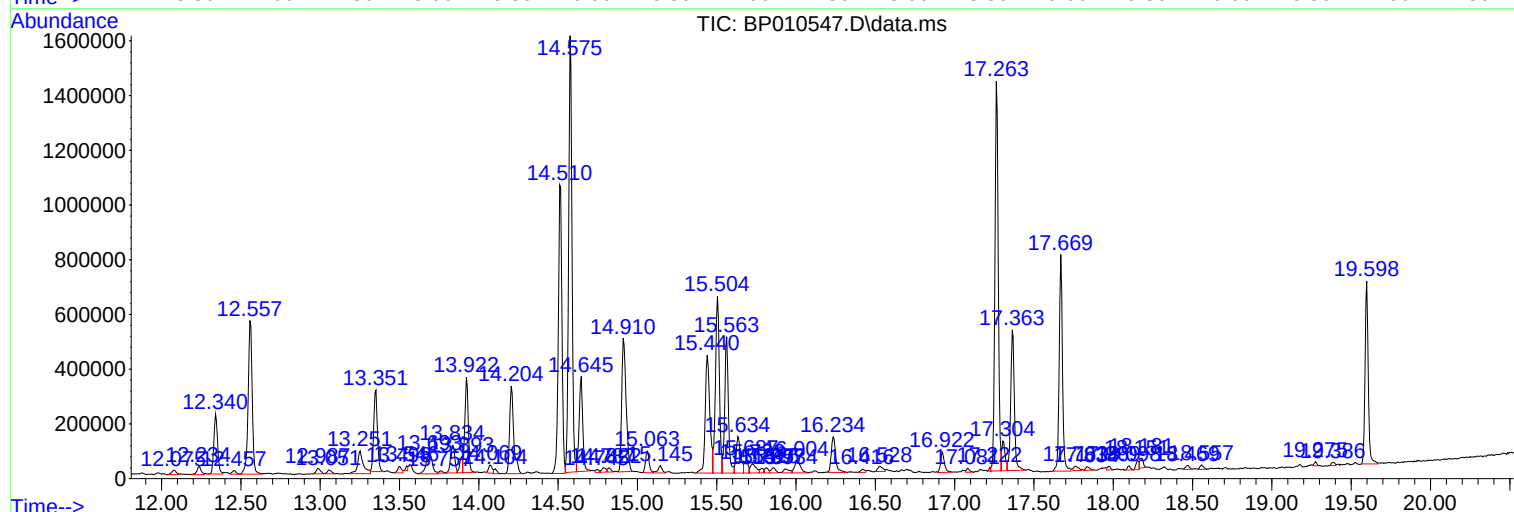
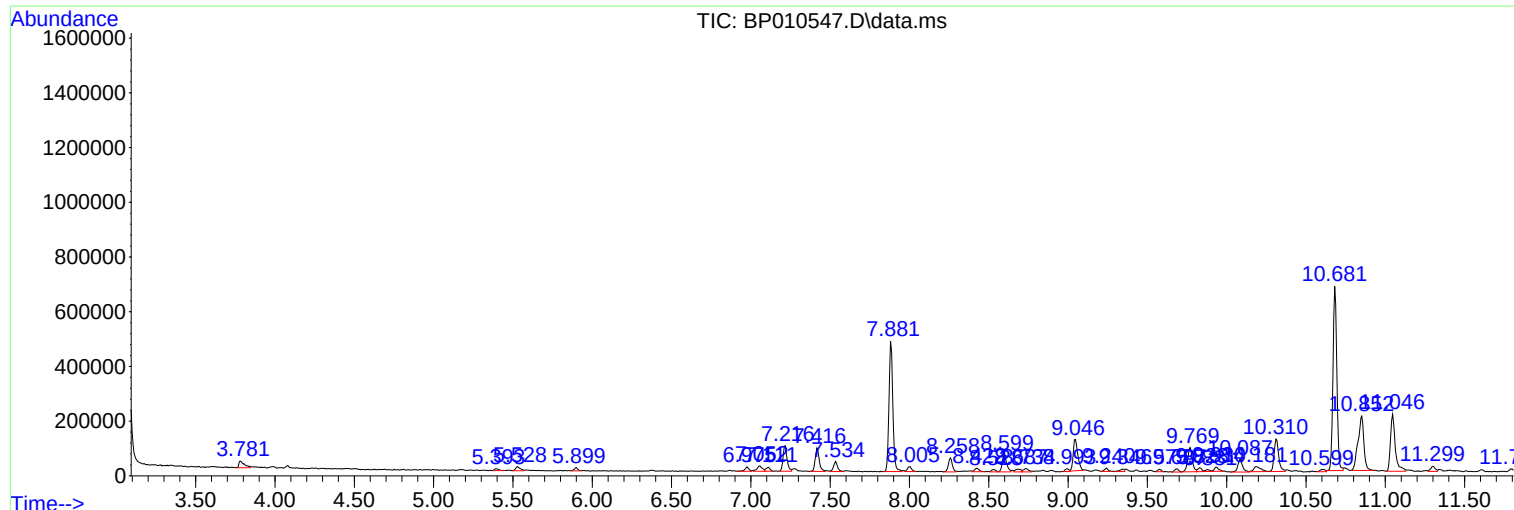
Sum of corrected areas: 27717325

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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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



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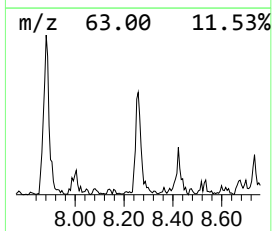
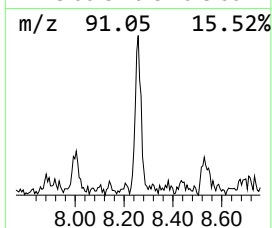
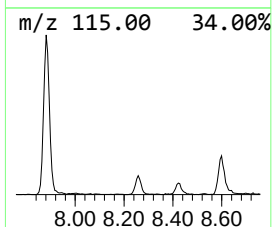
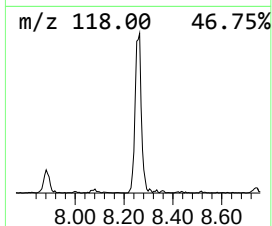
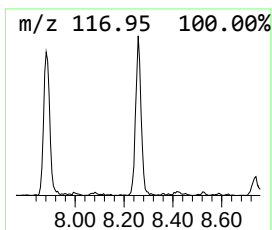
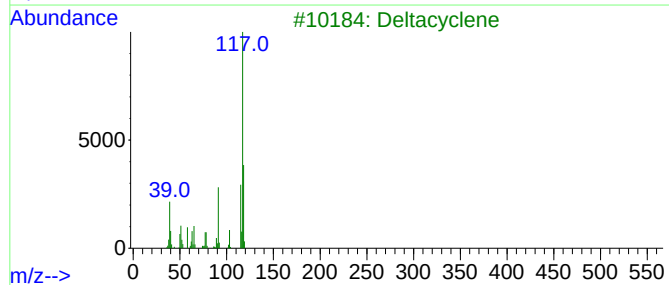
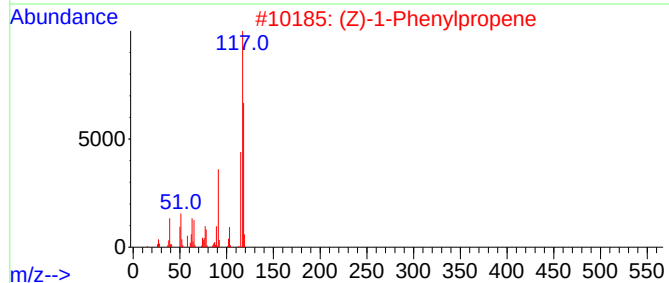
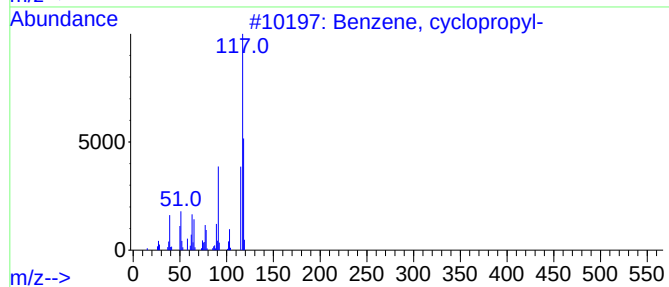
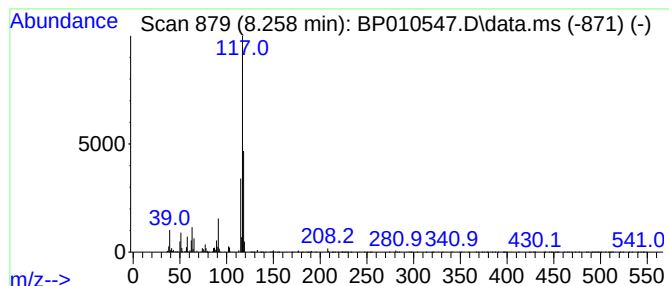
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Benzene, cyclopropyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.258	2.18 ng/ul	86994	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
2			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	74
3			Deltacyclene	118	C9H10	007785-10-6	64
4			Indane	118	C9H10	000496-11-7	64
5			Indane	118	C9H10	000496-11-7	60



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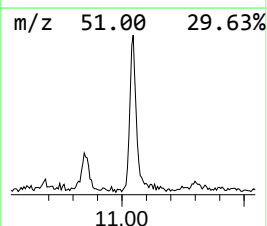
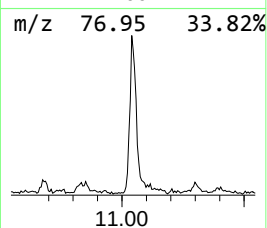
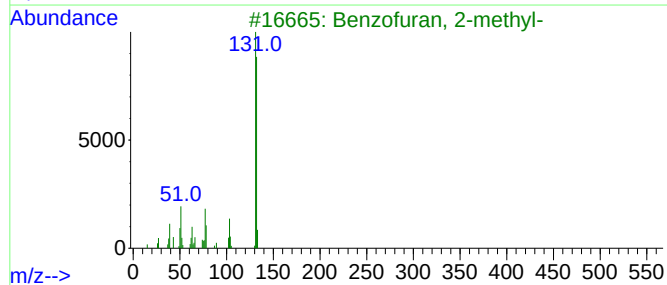
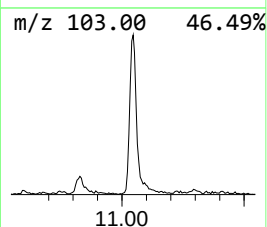
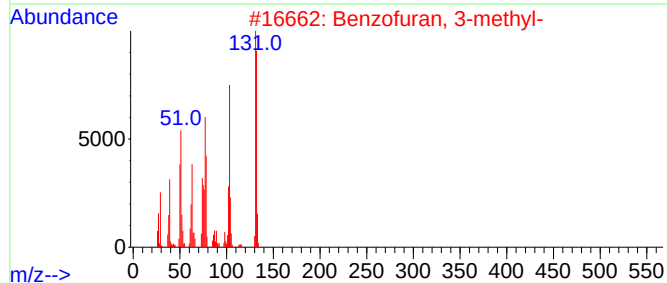
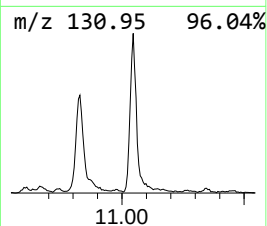
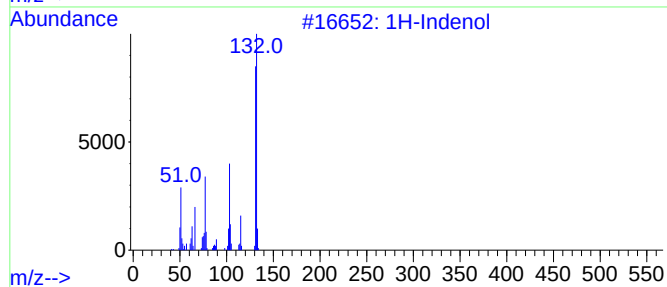
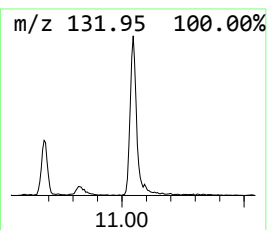
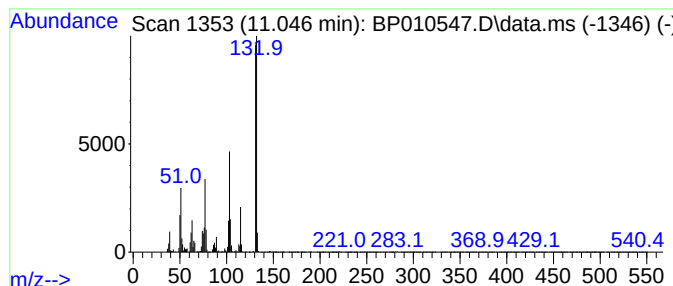
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1H-Indenol Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.046	7.54 ng/ul	430447	Naphthalene-d8	10.681

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indenol	132	C9H8O	056631-57-3	95
2		Benzofuran, 3-methyl-	132	C9H8O	021535-97-7	90
3		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
4		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	87
5		Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	83



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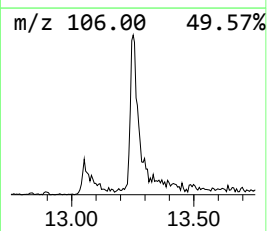
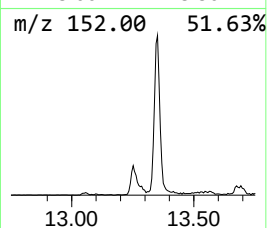
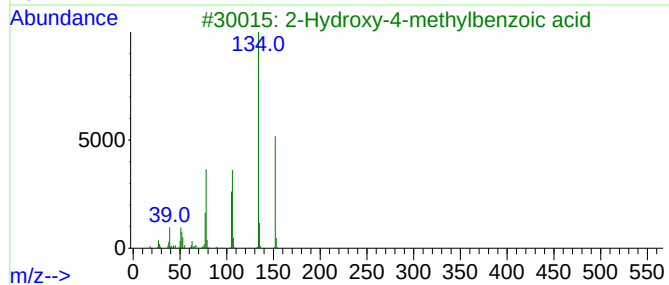
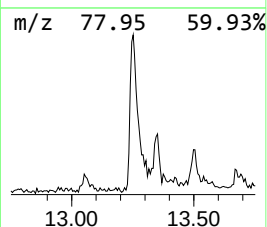
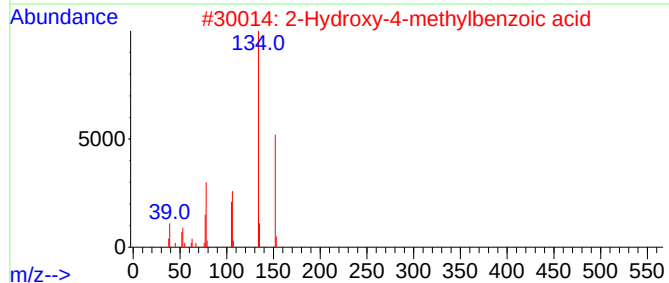
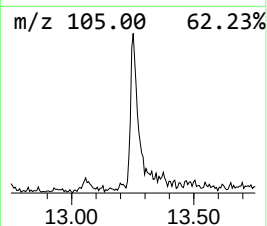
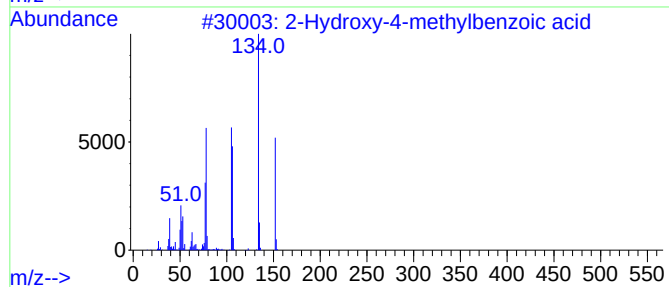
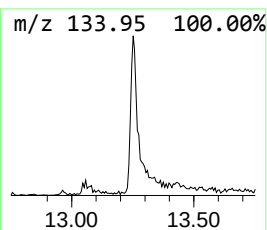
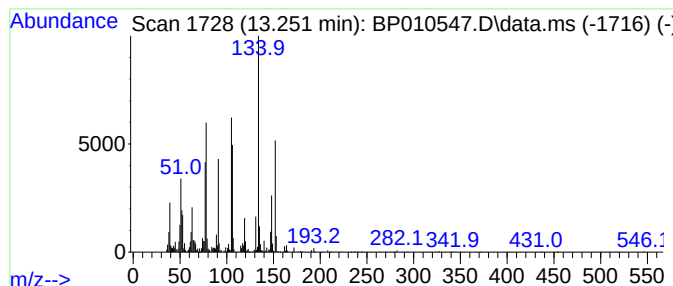
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 2-Hydroxy-4-methylbenzoic acid Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.251	2.51 ng/ul	210408	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hydroxy-4-methylbenzoic acid	152	C8H8O3	000050-85-1	95
2			2-Hydroxy-4-methylbenzoic acid	152	C8H8O3	000050-85-1	90
3			2-Hydroxy-4-methylbenzoic acid	152	C8H8O3	000050-85-1	76
4			5-Methylsalicylic acid	152	C8H8O3	000089-56-5	76
5			5-Methylsalicylic acid	152	C8H8O3	000089-56-5	60



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTD0DL

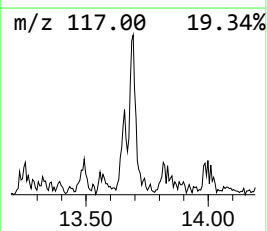
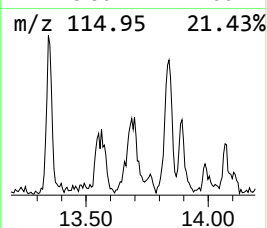
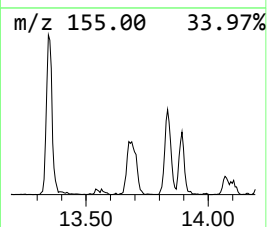
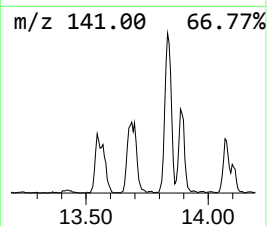
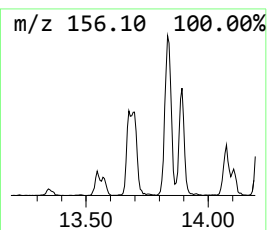
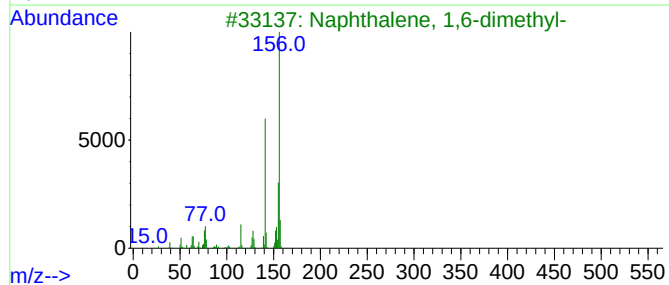
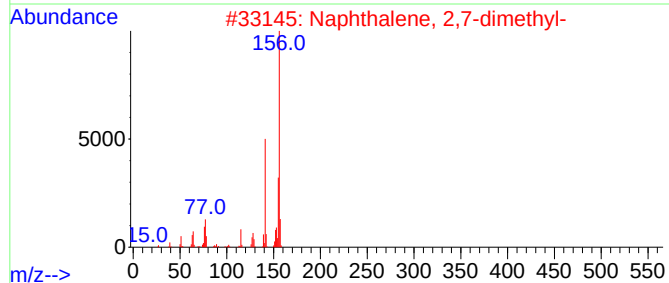
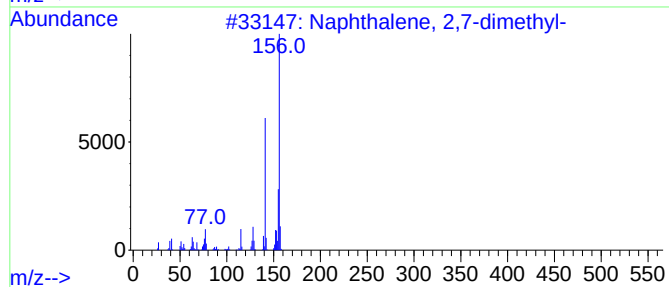
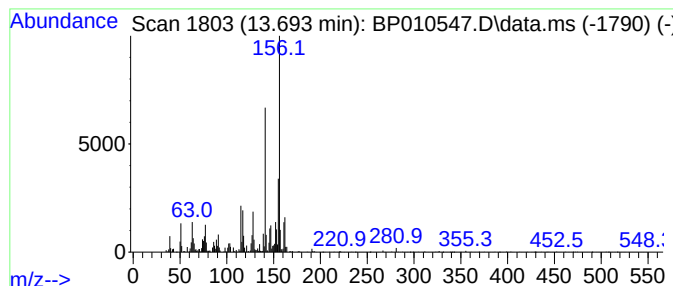
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2,7-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.693	2.39 ng/ul	200842	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	96
2			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	95
4			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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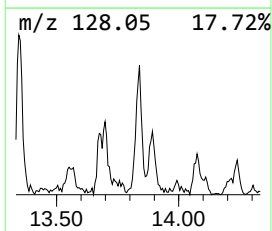
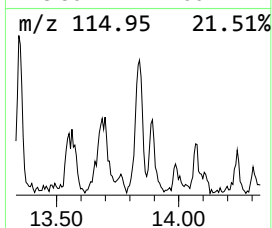
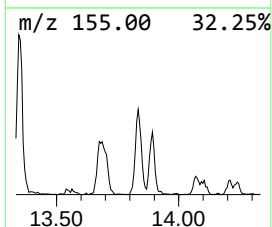
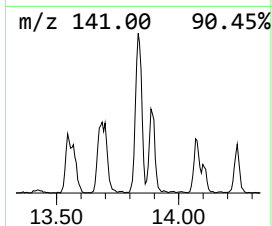
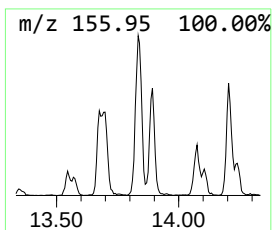
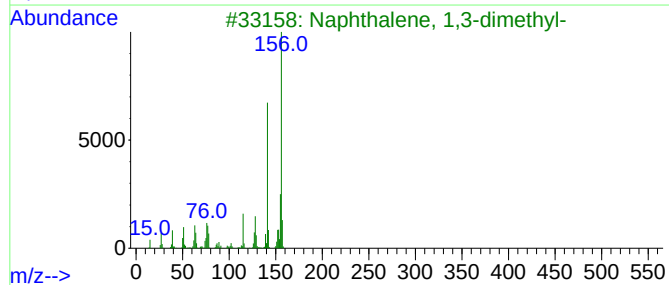
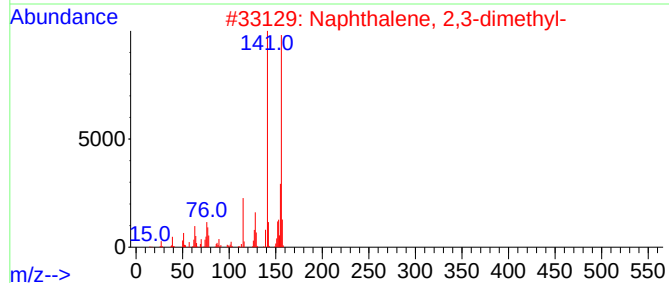
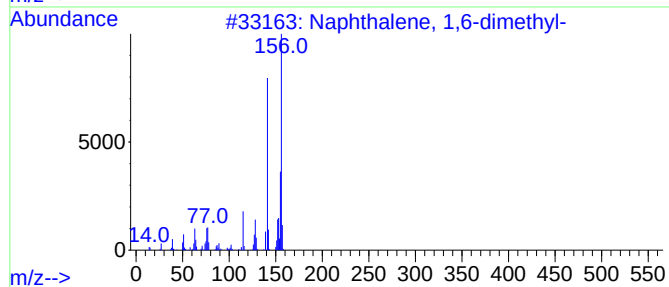
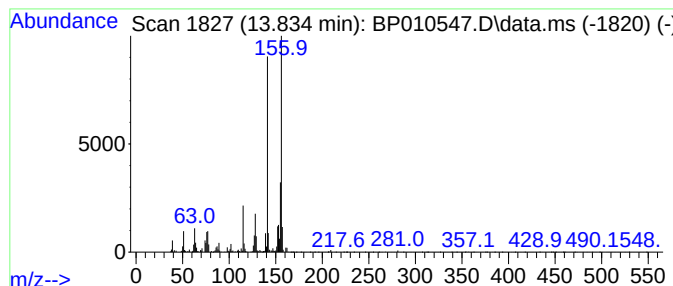
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 1,6-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.834	2.25 ng/ul	188872	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	98
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	98
3			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	97
4			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97
5			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	97



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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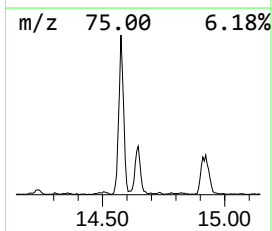
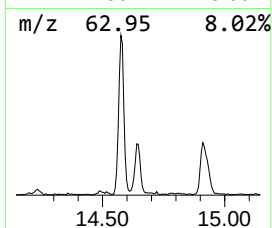
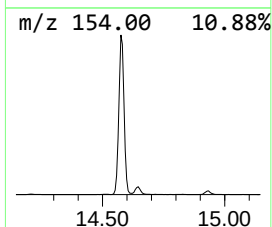
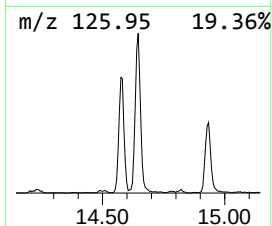
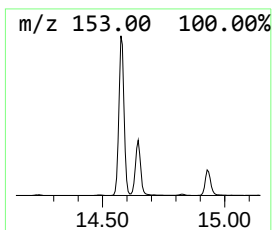
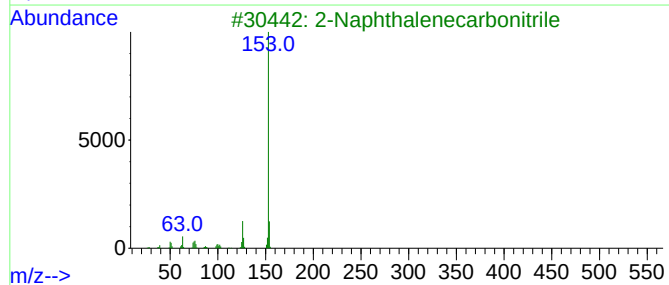
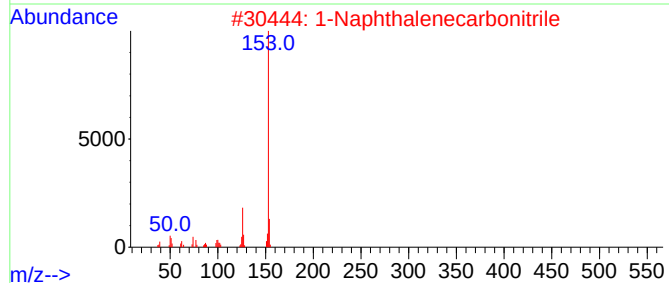
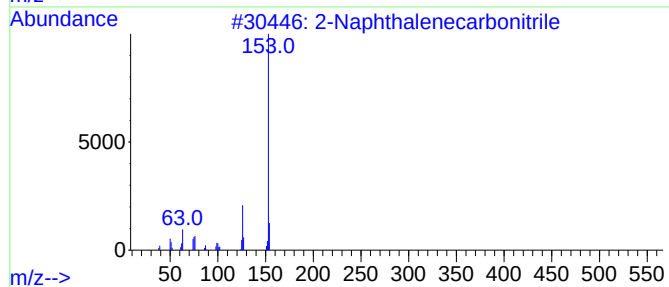
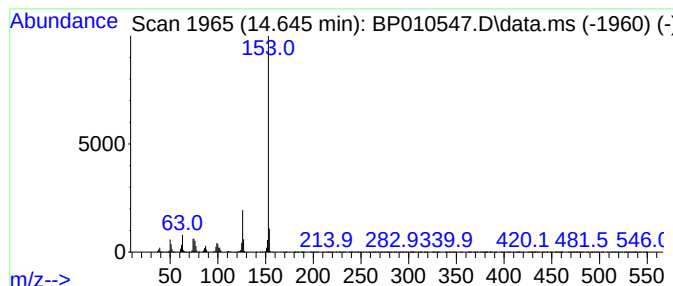
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Naphthalenecarbonitrile Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.646	6.42 ng/ul	539221	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	97		
2	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94		
3	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	94		
4	1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	94		
5	2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	91		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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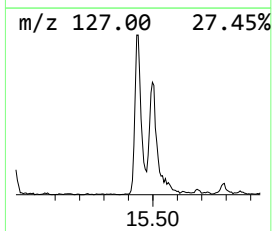
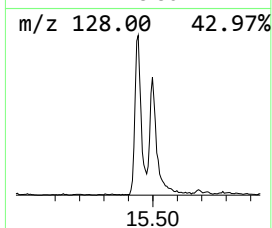
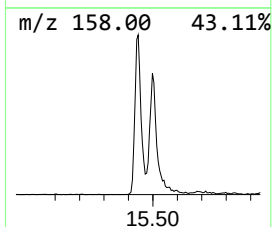
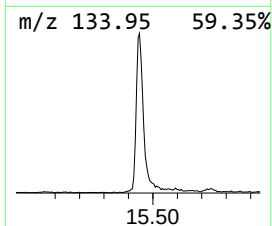
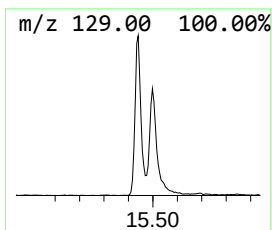
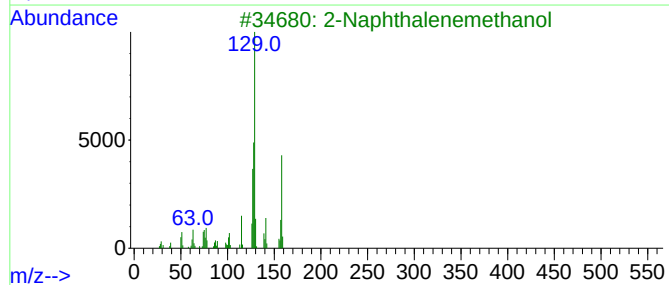
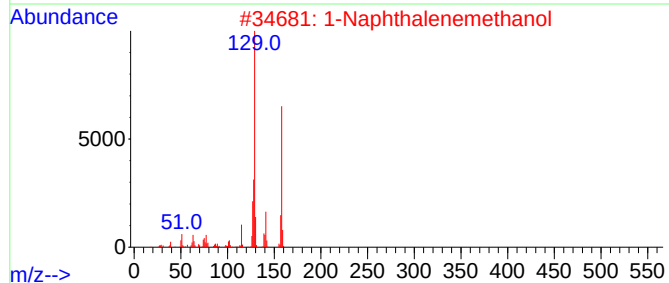
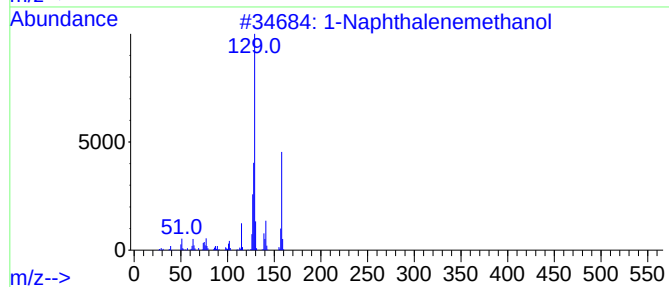
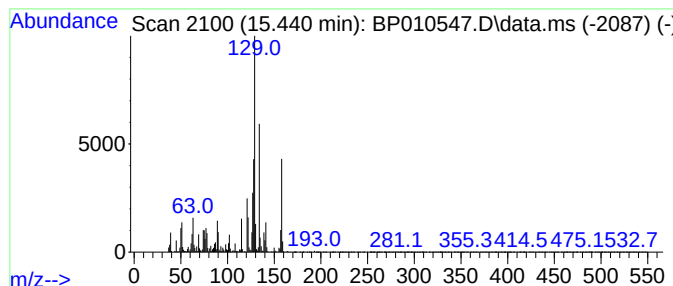
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Naphthalenemethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	10.34 ng/ul	868071	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Naphthalenemethanol			158	C11H10O	004780-79-4	98
2	1-Naphthalenemethanol			158	C11H10O	004780-79-4	94
3	2-Naphthalenemethanol			158	C11H10O	001592-38-7	93
4	1-Naphthalenemethanol			158	C11H10O	004780-79-4	90
5	1-Naphthalenemethanol			158	C11H10O	004780-79-4	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
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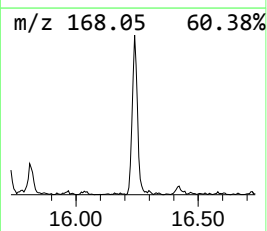
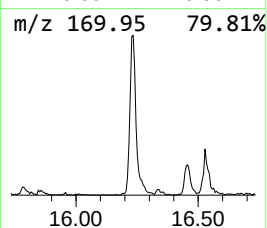
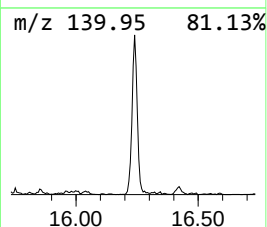
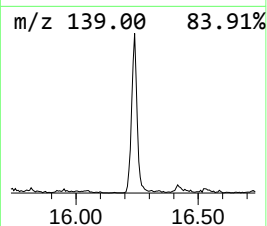
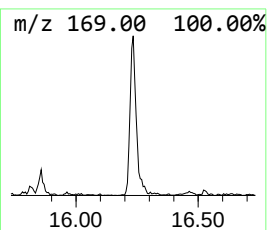
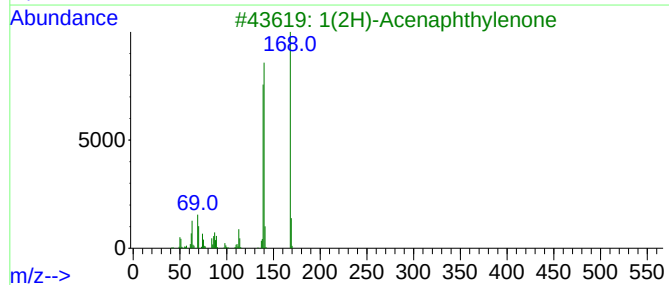
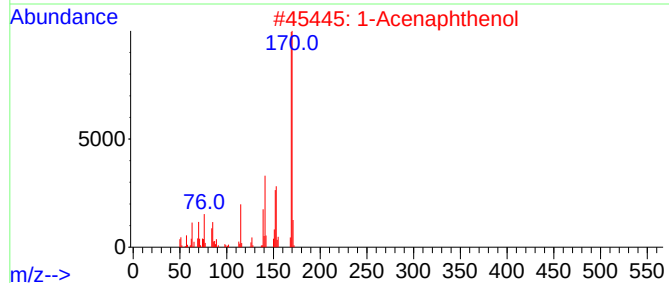
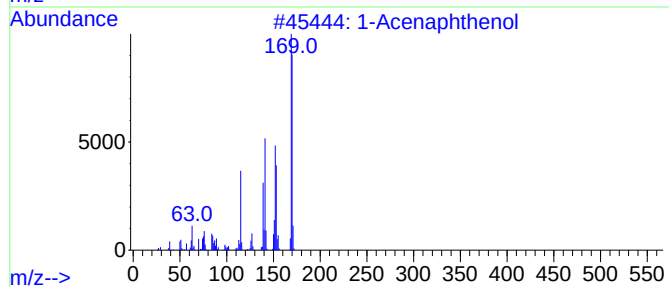
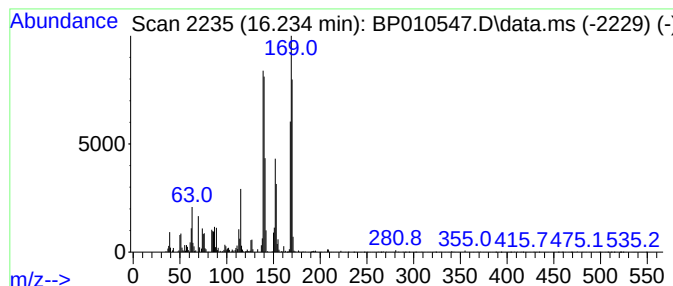
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Acenaphthenol Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.234	2.37 ng/ul	247371	Phenanthrene-d10	17.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Acenaphthenol	170	C12H10O	006306-07-6	89
2			1-Acenaphthenol	170	C12H10O	006306-07-6	87
3			1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	55
4			1-Acenaphthenol	170	C12H10O	006306-07-6	41
5			Benzaldehyde, 2-fluoro-5-hydroxy...	170	C8H7FO3	079418-76-1	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010547.D
Acq On : 25 May 2022 21:31
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 17 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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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
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1H-Indenol	11.046	7.5	ng/u1	430447	2	10.681	1141650	20.0
2-Hydroxy-4-met...	13.251	2.5	ng/u1	210408	3	14.516	1678920	20.0
Naphthalene, 2,...	13.693	2.4	ng/u1	200842	3	14.516	1678920	20.0
Naphthalene, 1,...	13.834	2.3	ng/u1	188872	3	14.516	1678920	20.0
2-Naphthaleneca...	14.646	6.4	ng/u1	539221	3	14.516	1678920	20.0
1-Naphthaleneme...	15.440	10.3	ng/u1	868071	3	14.516	1678920	20.0
1-Acenaphthenol	16.234	2.4	ng/u1	247371	4	17.263	2089870	20.0