

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
 Data File : BP010540.D
 Acq On : 25 May 2022 17:34
 Operator : CG/JU
 Sample : N2950-10DL 5X
 Misc :
 ALS Vial : 11 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: BP010540.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.776	115	117	125	rBV	28021	56010	1.99%	0.174%
2	5.528	410	415	425	rVV3	14461	30424	1.08%	0.094%
3	5.899	472	478	482	rBV2	12575	21845	0.78%	0.068%
4	6.975	654	661	667	rVV2	20052	35171	1.25%	0.109%
5	7.052	667	674	679	rVV3	24303	55350	1.96%	0.172%
6	7.111	680	684	693	rVV2	17184	31339	1.11%	0.097%
7	7.217	694	702	709	rVV	112526	202444	7.19%	0.628%
8	7.275	709	712	719	rVV5	11831	22653	0.80%	0.070%
9	7.417	731	736	747	rVB	98526	172650	6.13%	0.536%
10	7.534	749	756	763	rVV	41748	67449	2.39%	0.209%
11	7.881	808	815	827	rBV	432508	728919	25.87%	2.261%
12	7.999	831	835	840	rVB	20432	33634	1.19%	0.104%
13	8.258	874	879	885	rBV	59420	95925	3.41%	0.298%
14	8.422	902	907	919	rVB2	14922	29082	1.03%	0.090%
15	8.599	930	937	946	rBV2	72297	141305	5.02%	0.438%
16	8.728	956	959	967	rVB6	10878	21367	0.76%	0.066%
17	8.993	998	1004	1007	rBV3	13462	22765	0.81%	0.071%
18	9.046	1007	1013	1022	rVV	146590	301052	10.69%	0.934%
19	9.181	1030	1036	1041	rVV7	7643	17673	0.63%	0.055%
20	9.240	1041	1046	1053	rVB6	12024	22203	0.79%	0.069%
21	9.352	1058	1065	1066	rVV7	12467	24841	0.88%	0.077%
22	9.369	1066	1068	1073	rVV5	11860	18321	0.65%	0.057%
23	9.569	1099	1102	1109	rVB2	10401	18134	0.64%	0.056%
24	9.681	1116	1121	1127	rBV5	14519	23465	0.83%	0.073%
25	9.769	1127	1136	1143	rBV	113880	224111	7.96%	0.695%
26	9.828	1143	1146	1151	rVB2	13916	20485	0.73%	0.064%
27	9.934	1160	1164	1168	rBV2	16872	24725	0.88%	0.077%
28	10.081	1182	1189	1198	rBV2	52391	102245	3.63%	0.317%
29	10.187	1201	1207	1218	rBV3	21355	60594	2.15%	0.188%
30	10.310	1222	1228	1237	rBV2	152495	283417	10.06%	0.879%
31	10.605	1268	1278	1283	rBV10	10832	24607	0.87%	0.076%
32	10.681	1283	1291	1300	rVV	637336	1097502	38.96%	3.404%
33	10.746	1300	1302	1308	rVB6	12607	18728	0.66%	0.058%
34	10.852	1308	1320	1330	rBV2	222721	592091	21.02%	1.837%
35	11.046	1346	1353	1368	rBV	256489	523020	18.57%	1.622%
36	11.299	1390	1396	1401	rVV3	23756	42422	1.51%	0.132%

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

37	11.793	1474	1480	1486	rVV4	13056	22602	0.80%	0.070%
38	12.081	1522	1529	1537	rVB4	19892	42932	1.52%	0.133%
39	12.240	1547	1556	1560	rBV	32685	59682	2.12%	0.185%
40	12.340	1566	1573	1582	rBV	255382	431867	15.33%	1.340%

41	12.457	1589	1593	1599	rVB2	13250	24510	0.87%	0.076%
42	12.557	1599	1610	1622	rBV	697596	1188257	42.18%	3.686%
43	12.987	1676	1683	1689	rVV2	25079	53154	1.89%	0.165%
44	13.052	1689	1694	1701	rVB8	19089	41449	1.47%	0.129%
45	13.252	1722	1728	1738	rVV2	129766	282928	10.04%	0.878%

46	13.352	1738	1745	1754	rVB	384955	598599	21.25%	1.857%
47	13.499	1765	1770	1774	rBV2	25018	44197	1.57%	0.137%
48	13.563	1774	1781	1790	rVB4	35017	104239	3.70%	0.323%
49	13.693	1791	1803	1811	rBV4	82933	225377	8.00%	0.699%
50	13.834	1820	1827	1833	rBV2	122915	231050	8.20%	0.717%

51	13.893	1833	1837	1838	rVV	74111	91539	3.25%	0.284%
52	13.922	1838	1842	1851	rVB	433127	611158	21.69%	1.896%
53	14.069	1863	1867	1871	rVV	33901	54462	1.93%	0.169%
54	14.104	1871	1873	1878	rVB3	21127	25668	0.91%	0.080%
55	14.204	1884	1890	1901	rBV	385547	691124	24.53%	2.144%

56	14.363	1912	1917	1924	rVB8	10132	19816	0.70%	0.061%
57	14.516	1933	1943	1948	rBV	1024025	1623885	57.64%	5.037%
58	14.575	1948	1953	1959	rVV	1883579	2817106	100.00%	8.738%
59	14.646	1959	1965	1971	rVB	418090	649410	23.05%	2.014%
60	14.787	1986	1989	1992	rVV5	13688	20776	0.74%	0.064%

61	14.822	1992	1995	2001	rVB3	14431	23745	0.84%	0.074%
62	14.910	2004	2010	2021	rVB2	580563	1122210	39.84%	3.481%
63	15.063	2030	2036	2042	rBV	108563	159373	5.66%	0.494%
64	15.146	2045	2050	2055	rVB3	32428	47816	1.70%	0.148%
65	15.440	2085	2100	2106	rBV2	617601	1163002	41.28%	3.607%

66	15.504	2106	2111	2116	rVV2	795529	1292867	45.89%	4.010%
67	15.563	2116	2121	2128	rVV	597834	880742	31.26%	2.732%
68	15.634	2128	2133	2139	rVV2	158938	291347	10.34%	0.904%
69	15.687	2139	2142	2145	rVV2	42016	58731	2.08%	0.182%
70	15.728	2145	2149	2156	rVB7	27069	44271	1.57%	0.137%

71	15.816	2161	2164	2167	rVV4	19688	29563	1.05%	0.092%
72	15.857	2167	2171	2176	rVB2	22365	38119	1.35%	0.118%
73	15.934	2179	2184	2191	rBV4	20042	41924	1.49%	0.130%
74	16.004	2191	2196	2205	rVB4	54385	126314	4.48%	0.392%
75	16.234	2229	2235	2243	rBV2	162953	294812	10.47%	0.914%

76	16.416	2261	2266	2269	rBV5	17451	24835	0.88%	0.077%
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 Title : SVOA CALIBRATION

77	16.534	2280	2286	2290	rBV8	23469	55265	1.96%	0.171%
78	16.698	2311	2314	2323	rVB6	18699	36788	1.31%	0.114%
79	16.922	2347	2352	2362	rVB2	94795	158389	5.62%	0.491%
80	17.087	2376	2380	2387	rVB3	15420	24589	0.87%	0.076%
81	17.228	2400	2404	2405	rVV2	18840	22858	0.81%	0.071%
82	17.263	2405	2410	2415	rVV	1360359	2020313	71.72%	6.267%
83	17.310	2415	2418	2421	rVV	129868	188480	6.69%	0.585%
84	17.363	2422	2427	2438	rVB2	631946	990286	35.15%	3.072%
85	17.669	2470	2479	2490	rBV	988303	1467342	52.09%	4.551%
86	17.763	2490	2495	2503	rVV7	25397	58825	2.09%	0.182%
87	17.834	2503	2507	2514	rVV	19506	33960	1.21%	0.105%
88	17.939	2519	2525	2527	rVV6	13419	30095	1.07%	0.093%
89	17.975	2528	2531	2535	rVV3	16959	26253	0.93%	0.081%
90	18.098	2548	2552	2556	rBV	21712	32319	1.15%	0.100%
91	18.151	2556	2561	2563	rBV4	42033	57836	2.05%	0.179%
92	18.175	2563	2565	2571	rVB	44324	61687	2.19%	0.191%
93	18.469	2611	2615	2619	rVB4	20938	32125	1.14%	0.100%
94	18.557	2625	2630	2635	rBV2	20792	36333	1.29%	0.113%
95	19.275	2749	2752	2757	rVB	15721	22036	0.78%	0.068%
96	19.386	2767	2771	2778	rBV7	21220	36306	1.29%	0.113%
97	19.598	2801	2807	2820	rBV	868264	1221331	43.35%	3.788%
98	21.351	3099	3105	3119	rBV	1549850	2238148	79.45%	6.942%
99	23.610	3483	3489	3494	rBV	416036	827675	29.38%	2.567%
100	23.769	3509	3516	3525	rVB2	844395	1756001	62.33%	5.447%

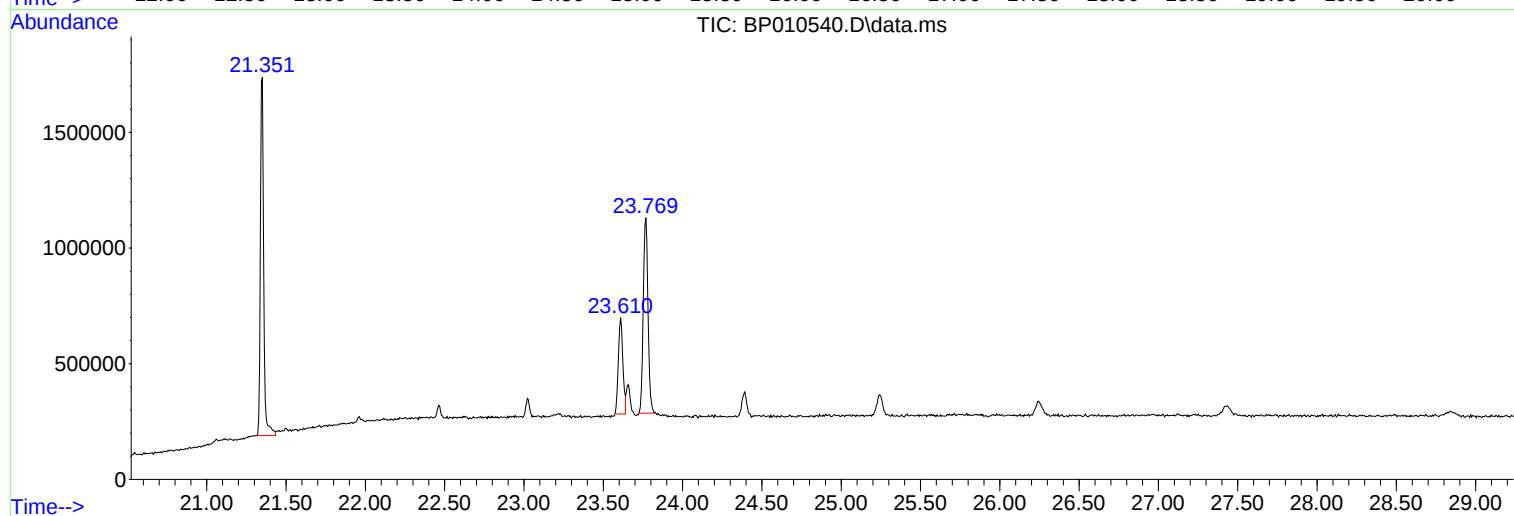
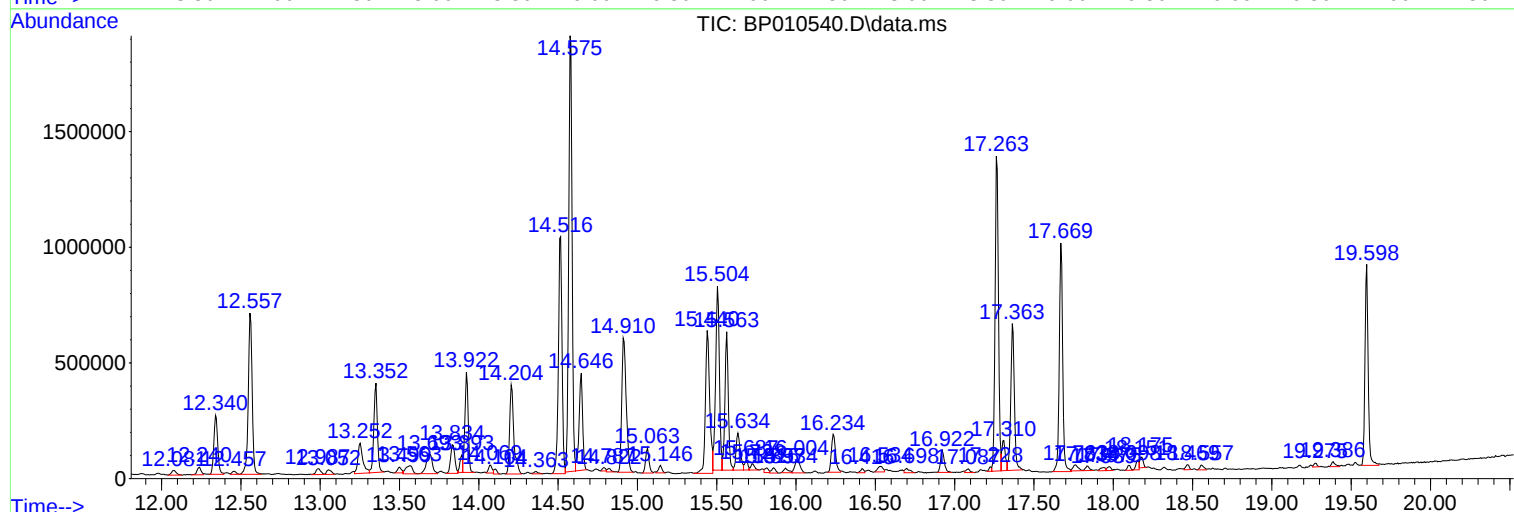
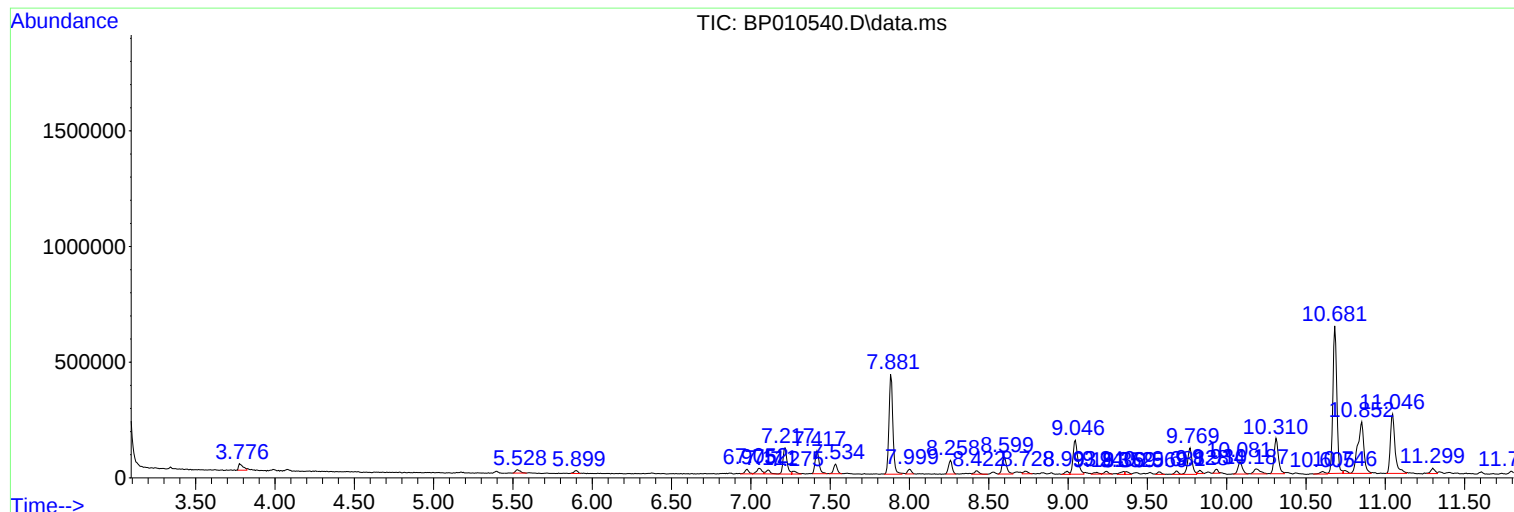
Sum of corrected areas: 32238666

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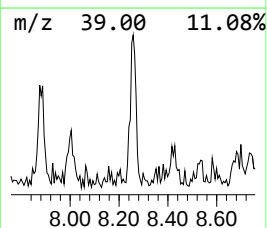
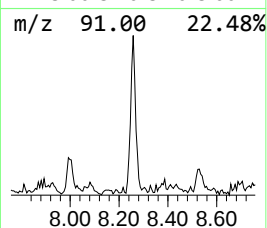
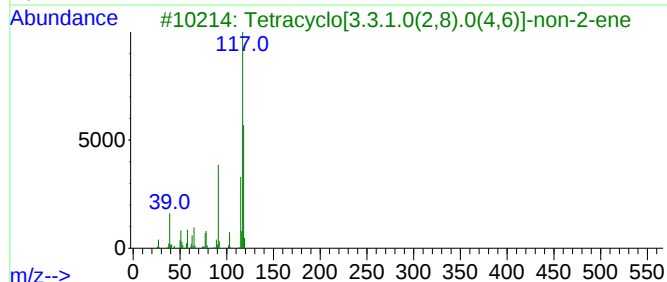
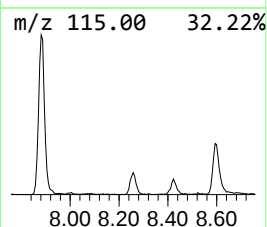
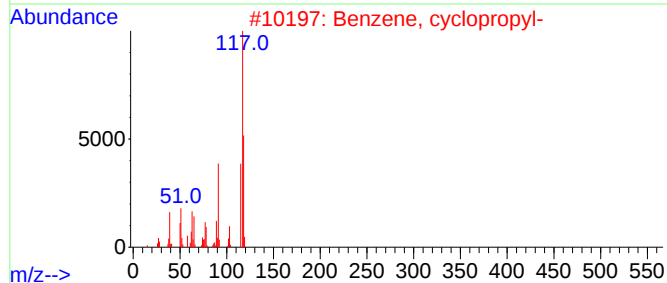
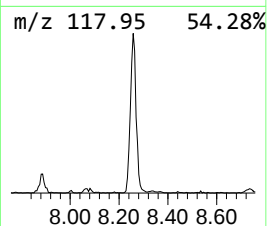
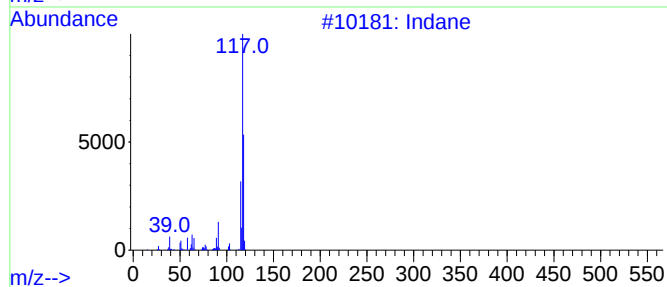
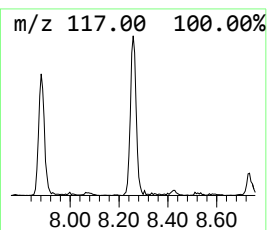
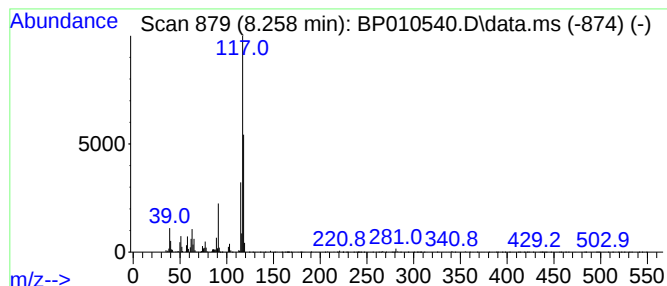
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.258	2.63 ng/ul	95925	1,4-Dichlorobenzene-d4	7.881

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	87
2			Benzene, cyclopropyl-	118	C9H10	000873-49-4	81
3			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
4			Deltacyclene	118	C9H10	007785-10-6	68
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	68



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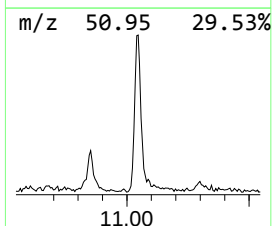
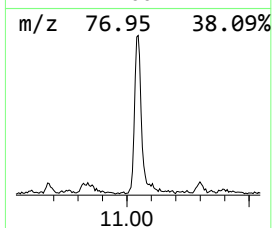
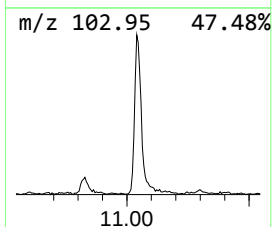
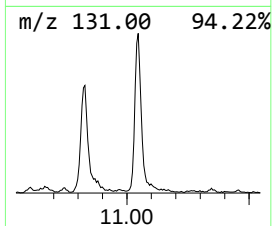
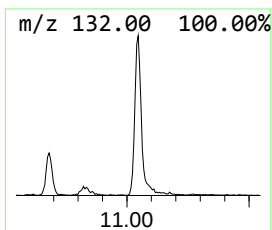
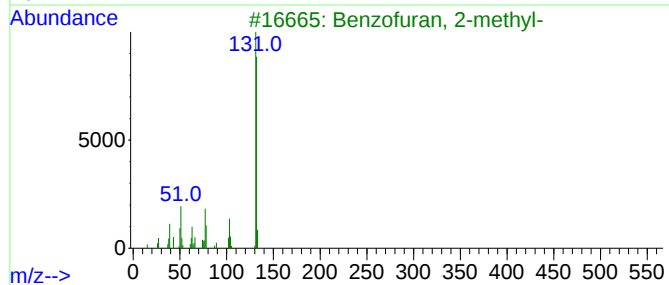
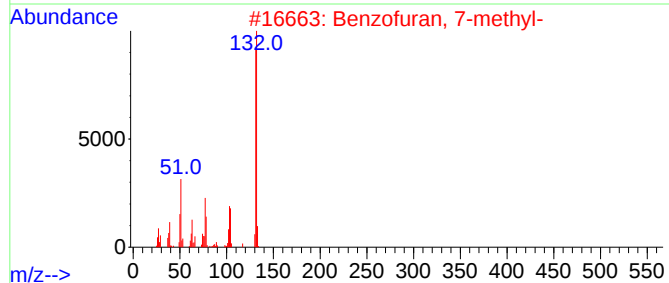
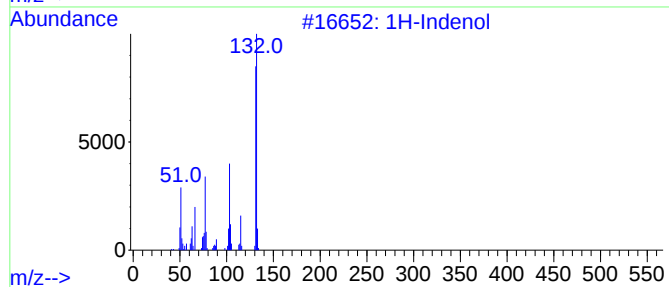
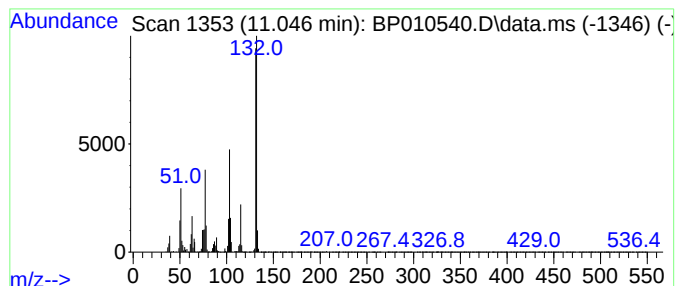
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051322.M
Quant Title : SVOA CALIBRATION

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	9.53 ng/ul	523020	Naphthalene-d8	10.681

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indenol	132	C9H8O	056631-57-3	94
2			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	86
3			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	83
4			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	81
5			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	72



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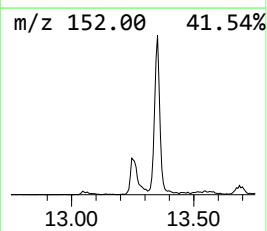
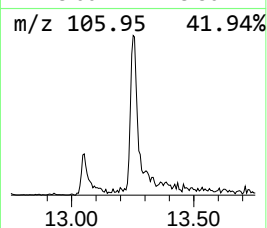
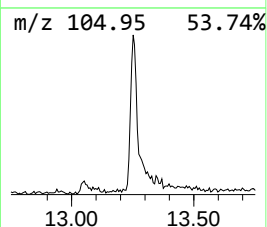
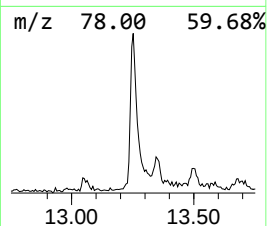
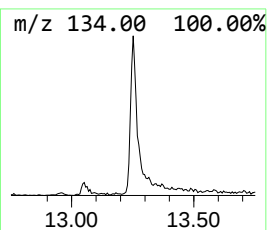
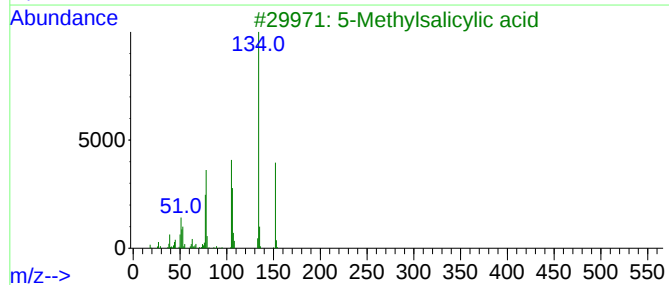
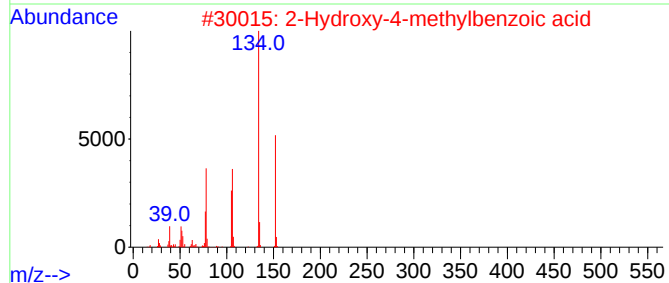
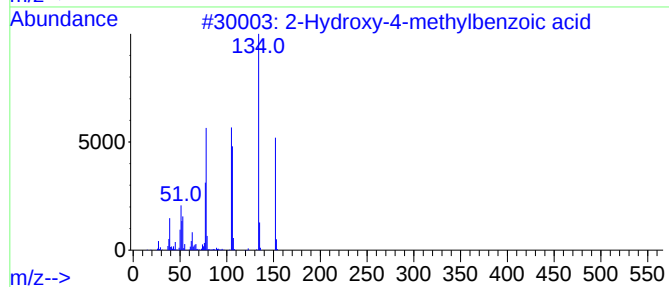
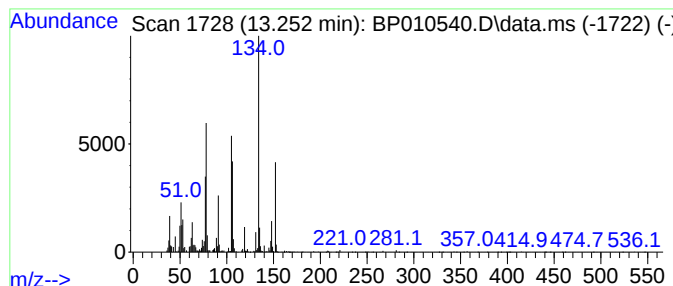
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Quant Title : SVOA CALIBRATION

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.252	3.48 ng/ul	282928	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			2-Hydroxy-4-methylbenzoic acid	152	C8H8O3	000050-85-1	96
2			2-Hydroxy-4-methylbenzoic acid	152	C8H8O3	000050-85-1	94
3			5-Methylsalicylic acid	152	C8H8O3	000089-56-5	93
4			5-Methylsalicylic acid	152	C8H8O3	000089-56-5	90
5			5-Methylsalicylic acid	152	C8H8O3	000089-56-5	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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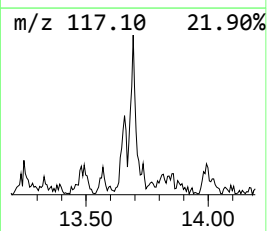
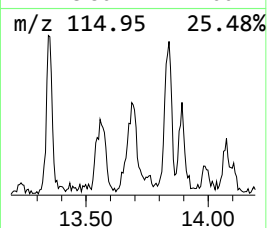
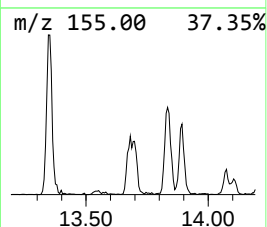
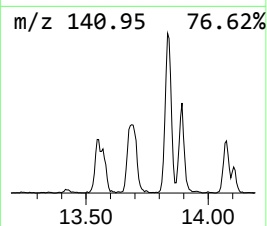
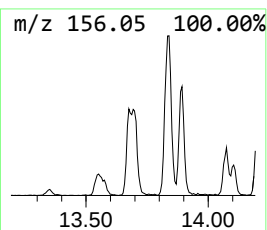
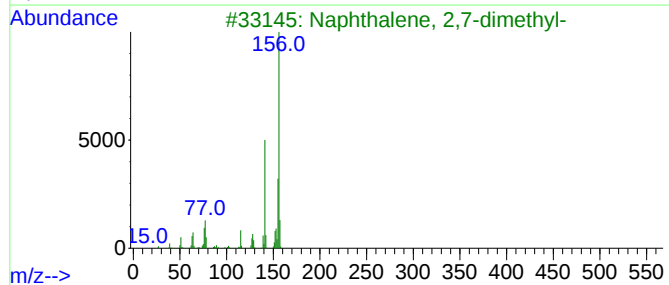
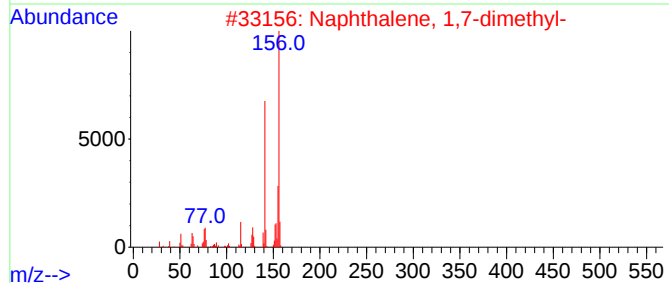
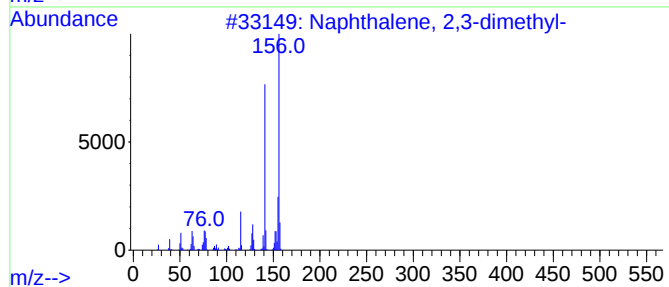
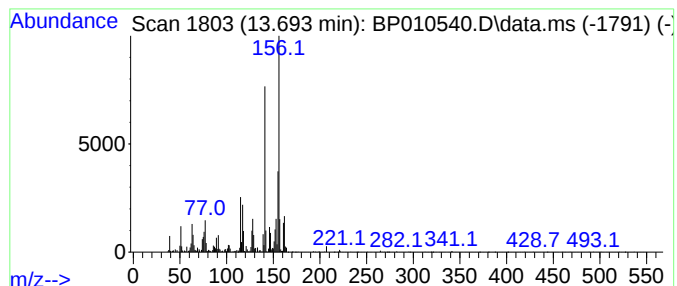
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Naphthalene, 2,3-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.693	2.78 ng/ul	225377	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	95
2			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	95
3			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	95
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	94
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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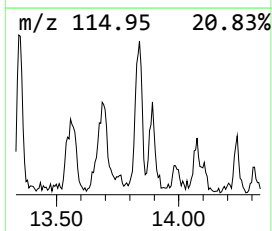
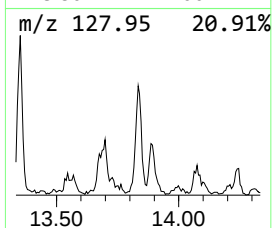
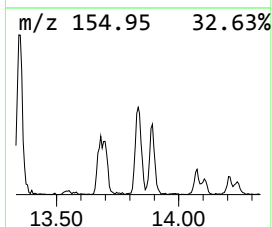
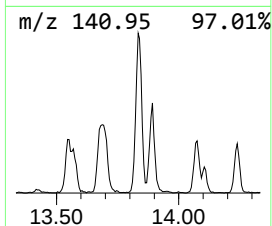
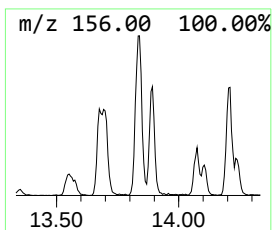
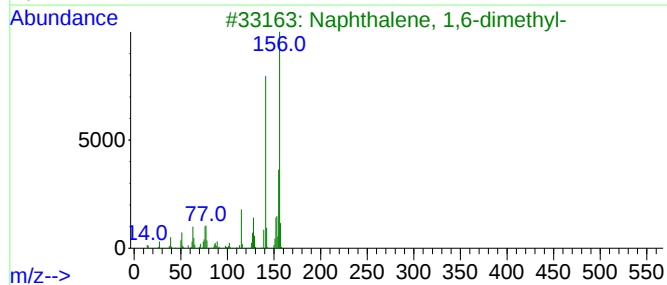
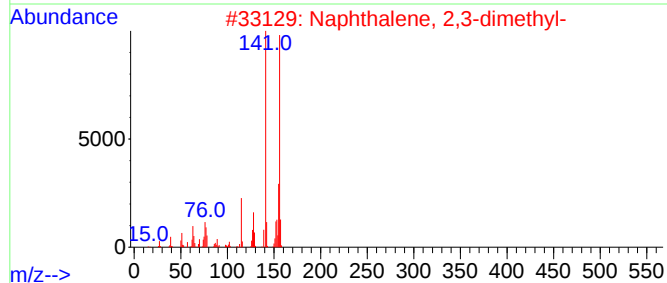
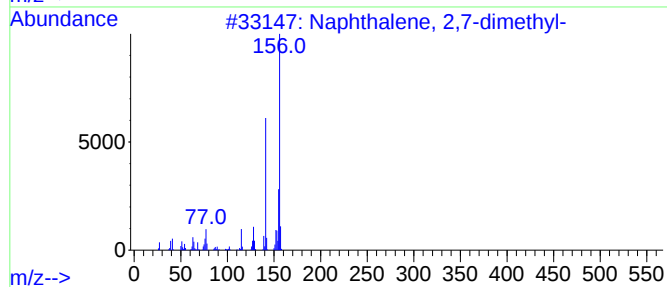
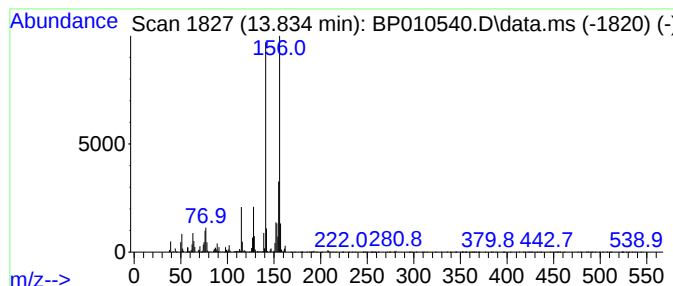
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Naphthalene, 2,7-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.834	2.85 ng/ul	231050	Acenaphthene-d10	14.516

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 2,7-dimethyl-	156	C12H12	000582-16-1	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 1,6-dimethyl-	156	C12H12	000575-43-9	96
4			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96
5			Naphthalene, 1,3-dimethyl-	156	C12H12	000575-41-7	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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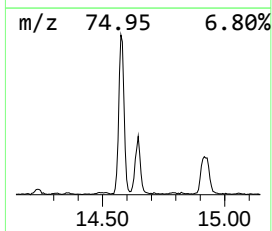
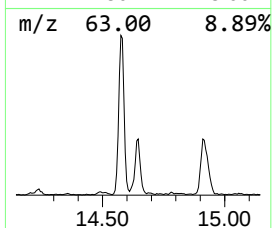
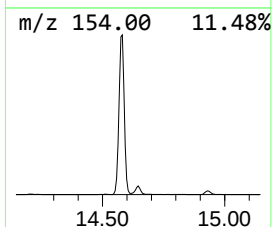
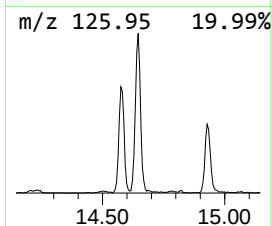
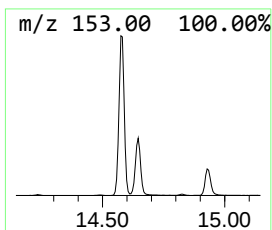
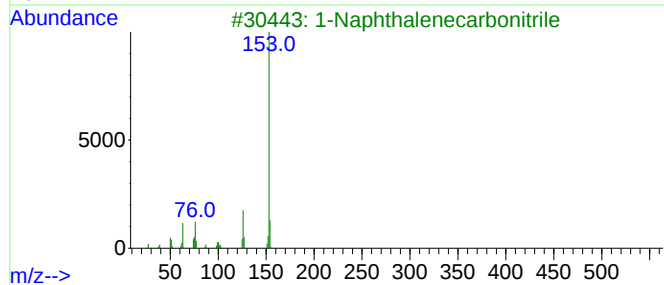
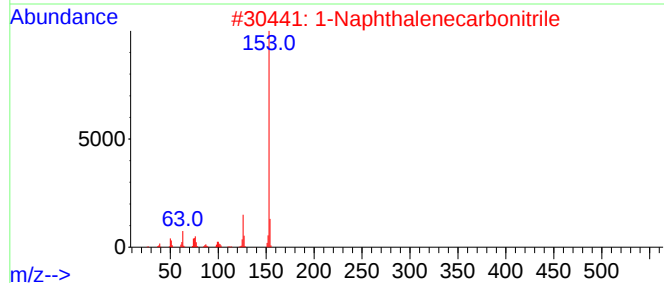
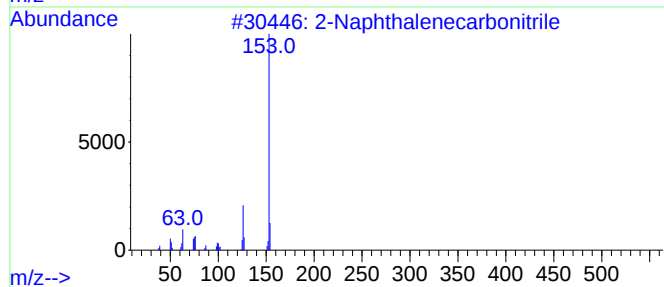
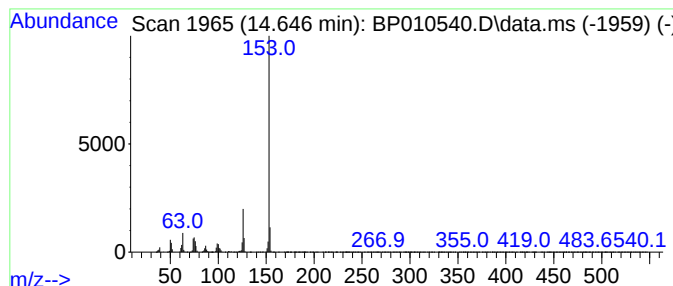
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 2-Naphthalenecarbonitrile Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.646	8.00 ng/ul	649410	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		1-Naphthalenecarbonitrile	153	C11H7N	000086-53-3	91
4		2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	90
5		2-Naphthalenecarbonitrile	153	C11H7N	000613-46-7	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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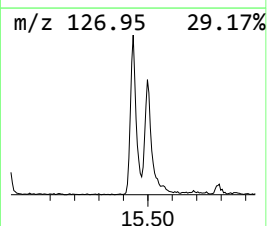
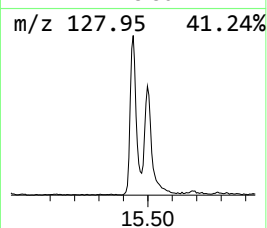
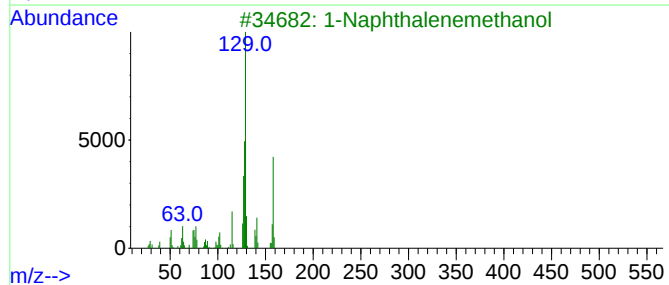
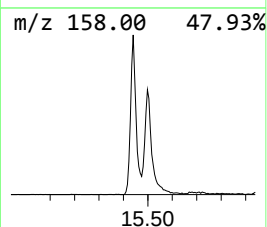
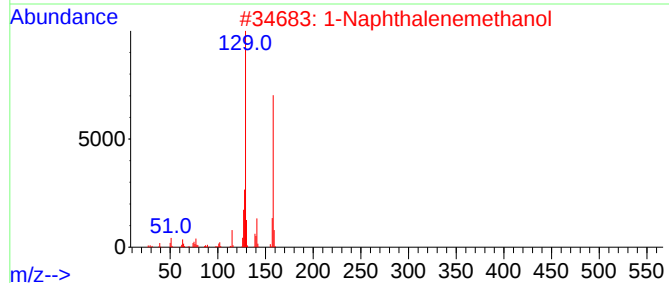
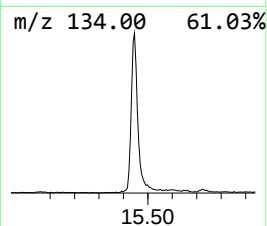
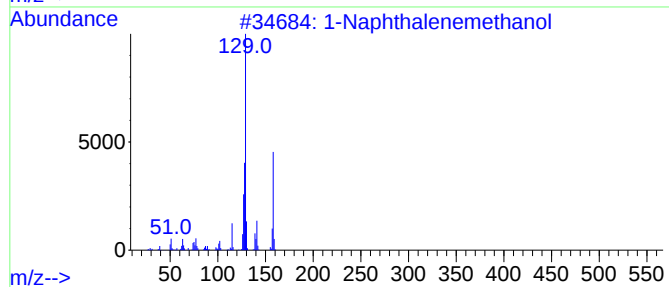
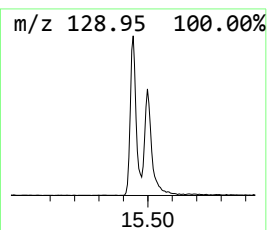
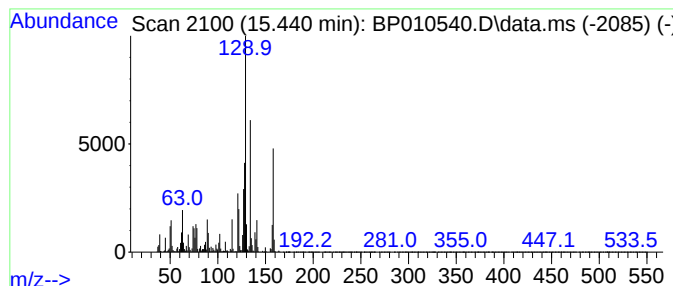
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 1-Naphthalenemethanol Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.440	14.32 ng/ul	1163000	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	91		
3	1-Naphthalenemethanol	158	C11H10O	004780-79-4	91		
4	1-Naphthalenemethanol	158	C11H10O	004780-79-4	91		
5	2-Naphthalenemethanol	158	C11H10O	001592-38-7	83		



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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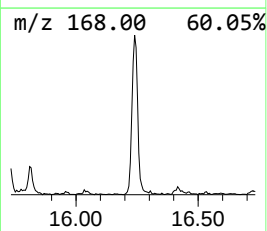
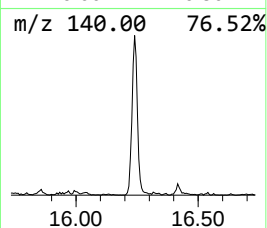
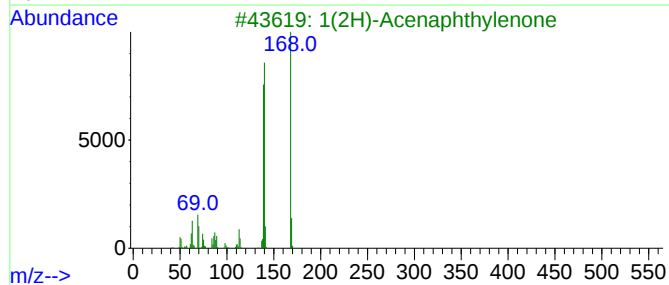
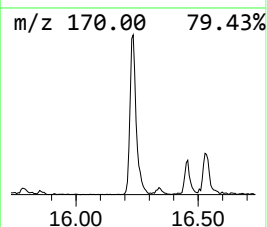
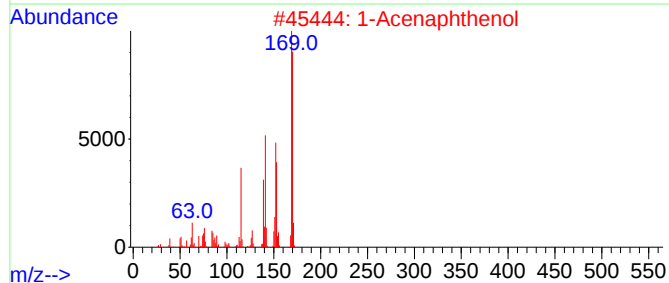
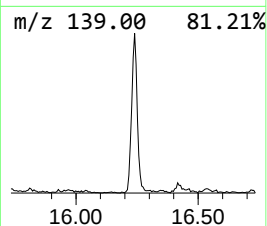
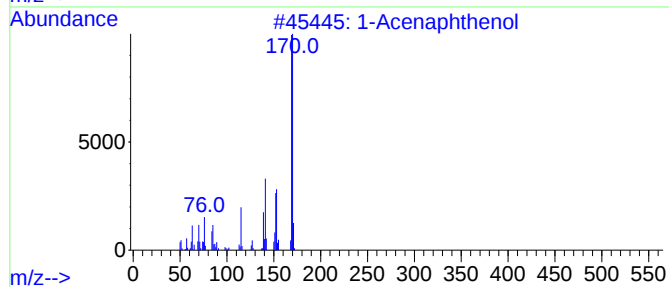
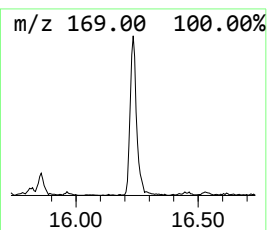
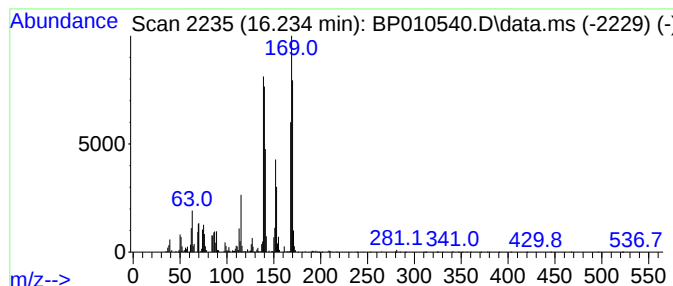
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 1-Acenaphthenol Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.234	2.92 ng/ul	294812	Phenanthrene-d10	17.263

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Acenaphthenol	170	C12H10O	006306-07-6	95
2			1-Acenaphthenol	170	C12H10O	006306-07-6	95
3			1(2H)-Acenaphthylenone	168	C12H8O	002235-15-6	83
4			1-Acenaphthenol	170	C12H10O	006306-07-6	46
5			1-Acenaphthenol, 3-methylamino-	199	C13H13NO	1000129-22-6	43



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP052522\
Data File : BP010540.D
Acq On : 25 May 2022 17:34
Operator : CG/JU
Sample : N2950-10DL 5X
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
DBTD0DL

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

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

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					#	RT	Resp	Conc
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1H-Indenol	11.046	9.5	ng/ul	523020	2	10.681	1097500	20.0
2-Hydroxy-4-met...	13.252	3.5	ng/ul	282928	3	14.516	1623890	20.0
Naphthalene, 2,...	13.693	2.8	ng/ul	225377	3	14.516	1623890	20.0
Naphthalene, 2,...	13.834	2.9	ng/ul	231050	3	14.516	1623890	20.0
2-Naphthaleneca...	14.646	8.0	ng/ul	649410	3	14.516	1623890	20.0
1-Naphthaleneme...	15.440	14.3	ng/ul	1163000	3	14.516	1623890	20.0
1-Acenaphthenol	16.234	2.9	ng/ul	294812	4	17.263	2020310	20.0