

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
 Data File : BP010338.D
 Acq On : 12 May 2022 23:38
 Operator : CG/JU
 Sample : N2801-02
 Misc :
 ALS Vial : 62 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 EXLB5

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-BP051022.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010338.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.311	34	38	41	rBV	24123	43303	0.84%	0.090%
2	3.346	41	44	53	rVV2	36855	67129	1.30%	0.140%
3	3.446	57	61	70	rVB	38131	62468	1.21%	0.131%
4	3.776	112	117	135	rBV2	109892	277426	5.38%	0.580%
5	3.999	152	155	161	rVB7	8019	11772	0.23%	0.025%
6	4.093	166	171	173	rBV2	10985	16475	0.32%	0.034%
7	4.128	173	177	186	rVB2	41846	64719	1.26%	0.135%
8	4.623	250	261	276	rVB2	103317	210015	4.07%	0.439%
9	4.887	303	306	311	rVB6	8735	15841	0.31%	0.033%
10	4.981	317	322	325	rBV2	10931	18640	0.36%	0.039%
11	5.175	350	355	362	rVB6	10168	17293	0.34%	0.036%
12	5.570	413	422	429	rBV	49181	87660	1.70%	0.183%
13	5.822	461	465	470	rVV2	12908	20550	0.40%	0.043%
14	5.887	470	476	477	rVV7	6000	11101	0.22%	0.023%
15	5.917	477	481	483	rVV2	11920	19700	0.38%	0.041%
16	5.952	483	487	496	rVB2	28249	57230	1.11%	0.120%
17	6.152	514	521	528	rBV2	8623	16967	0.33%	0.035%
18	6.393	555	562	570	rBV2	114706	193307	3.75%	0.404%
19	7.064	669	676	685	rBV	111122	214189	4.16%	0.447%
20	7.228	698	704	710	rBV	475238	790362	15.33%	1.651%
21	7.293	710	715	727	rVV	493938	789194	15.31%	1.649%
22	7.434	733	739	753	rVV	458019	810867	15.73%	1.694%
23	7.552	753	759	765	rVB	64478	104304	2.02%	0.218%
24	7.775	792	797	799	rBV2	8726	12978	0.25%	0.027%
25	7.805	799	802	807	rVB	16766	24088	0.47%	0.050%
26	7.905	809	819	833	rBV	443821	766010	14.86%	1.600%
27	8.022	833	839	848	rVB	38550	64138	1.24%	0.134%
28	8.275	876	882	889	rVB	23638	38786	0.75%	0.081%
29	8.399	896	903	906	rBV3	21384	35335	0.69%	0.074%
30	8.446	906	911	918	rVB3	21857	41913	0.81%	0.088%
31	8.558	923	930	933	rBV4	33405	58116	1.13%	0.121%
32	8.611	933	939	950	rVV	378040	651436	12.64%	1.361%
33	8.711	950	956	963	rVB2	57952	95960	1.86%	0.200%
34	8.864	975	982	987	rVB	20118	34951	0.68%	0.073%
35	9.016	998	1008	1010	rBV	63301	108303	2.10%	0.226%
36	9.064	1010	1016	1020	rBV	605109	1170171	22.70%	2.445%

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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-BP051022.M
 Title : SVOA CALIBRATION

37	9.228	1037	1044	1058	rVB	21477	54305	1.05%	0.113%
38	9.358	1058	1066	1072	rVB7	7494	17286	0.34%	0.036%
39	9.534	1086	1096	1103	rBV	155398	244481	4.74%	0.511%
40	9.711	1121	1126	1131	rVB6	13544	21301	0.41%	0.045%
41	9.787	1131	1139	1154	rBV	503366	901243	17.48%	1.883%
42	9.905	1154	1159	1163	rVV8	5561	14377	0.28%	0.030%
43	9.952	1163	1167	1179	rVB3	15907	36191	0.70%	0.076%
44	10.111	1185	1194	1203	rBV3	162525	310070	6.02%	0.648%
45	10.205	1203	1210	1214	rBV	32535	58878	1.14%	0.123%
46	10.246	1214	1217	1223	rVV4	12757	19209	0.37%	0.040%
47	10.328	1223	1231	1241	rVV	694493	1228695	23.84%	2.567%
48	10.410	1241	1245	1252	rVB6	16292	30083	0.58%	0.063%
49	10.705	1284	1295	1301	rBV	662299	1140555	22.13%	2.383%
50	10.758	1301	1304	1310	rVV2	42966	68081	1.32%	0.142%
51	10.840	1310	1318	1338	rVB	639529	1255306	24.35%	2.622%
52	11.146	1364	1370	1377	rVV5	9722	20498	0.40%	0.043%
53	11.610	1442	1449	1456	rBV6	7050	16926	0.33%	0.035%
54	11.810	1478	1483	1492	rBV8	8294	17508	0.34%	0.037%
55	12.075	1524	1528	1537	rVB4	19123	43588	0.85%	0.091%
56	12.187	1543	1547	1555	rBV9	7409	14239	0.28%	0.030%
57	12.269	1555	1561	1562	rBV5	9263	16128	0.31%	0.034%
58	12.293	1562	1565	1569	rVB3	10819	15865	0.31%	0.033%
59	12.334	1569	1572	1576	rBV4	6671	11251	0.22%	0.024%
60	12.428	1582	1588	1594	rVB5	9954	18096	0.35%	0.038%
61	12.522	1598	1604	1609	rVV2	19775	37428	0.73%	0.078%
62	12.581	1609	1614	1624	rVB4	34709	60344	1.17%	0.126%
63	12.999	1679	1685	1694	rVB	134449	225700	4.38%	0.472%
64	13.157	1707	1712	1717	rVB7	5781	10293	0.20%	0.022%
65	13.263	1723	1730	1738	rVV4	16346	35215	0.68%	0.074%
66	13.446	1758	1761	1765	rVV3	8605	11381	0.22%	0.024%
67	13.499	1765	1770	1779	rVV	307169	467079	9.06%	0.976%
68	13.569	1781	1782	1790	rVB6	8645	10417	0.20%	0.022%
69	13.663	1792	1798	1810	rBV	126427	232257	4.51%	0.485%
70	13.769	1812	1816	1820	rBV3	10348	13849	0.27%	0.029%
71	13.904	1829	1839	1840	rBV	82334	104957	2.04%	0.219%
72	13.940	1840	1845	1857	rVB	1731039	2439886	47.34%	5.097%
73	14.228	1885	1894	1904	rBV	1661453	2597122	50.39%	5.426%
74	14.304	1904	1907	1911	rVV5	11827	22161	0.43%	0.046%
75	14.351	1911	1915	1922	rVB2	24950	41711	0.81%	0.087%
76	14.457	1927	1933	1938	rBV7	6238	12702	0.25%	0.027%

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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

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77	14.534	1939	1946	1961	rVB	1124488	1693699	32.86%	3.538%
78	14.640	1961	1964	1971	rVB9	5826	10106	0.20%	0.021%
79	14.734	1971	1980	1997	rBV2	89371	245053	4.75%	0.512%
80	15.010	2022	2027	2032	rVB4	9677	18270	0.35%	0.038%
81	15.163	2046	2053	2058	rBV6	17945	32923	0.64%	0.069%
82	15.363	2081	2087	2095	rVB10	13443	28606	0.55%	0.060%
83	15.528	2105	2115	2128	rBV	2350314	3512996	68.15%	7.339%
84	15.646	2128	2135	2150	rVV	902111	1319118	25.59%	2.756%
85	15.769	2151	2156	2162	rVB	29022	46359	0.90%	0.097%
86	16.040	2198	2202	2209	rBV10	5500	11317	0.22%	0.024%
87	16.310	2244	2248	2253	rVB5	9266	13311	0.26%	0.028%
88	17.281	2406	2413	2424	rBV	1408701	2082381	40.40%	4.350%
89	17.381	2424	2430	2447	rVB	2792847	4212402	81.72%	8.800%
90	18.169	2559	2564	2571	rBV4	35714	60382	1.17%	0.126%
91	18.234	2573	2575	2581	rVB4	6549	10325	0.20%	0.022%
92	19.192	2734	2738	2744	rBV2	40461	55902	1.08%	0.117%
93	19.263	2746	2750	2754	rBV	35637	46557	0.90%	0.097%
94	19.398	2770	2773	2778	rBV7	11756	18423	0.36%	0.038%
95	19.616	2804	2810	2821	rBV	3971610	5154427	100.00%	10.768%
96	21.075	3054	3058	3065	rBV	105152	168335	3.27%	0.352%
97	21.363	3102	3107	3121	rBV	1754679	2403563	46.63%	5.021%
98	22.480	3292	3297	3306	rBV2	337043	529521	10.27%	1.106%
99	23.639	3486	3494	3501	rBV2	2408227	5007323	97.15%	10.461%
100	23.792	3513	3520	3529	rVB2	1131580	2266177	43.97%	4.734%

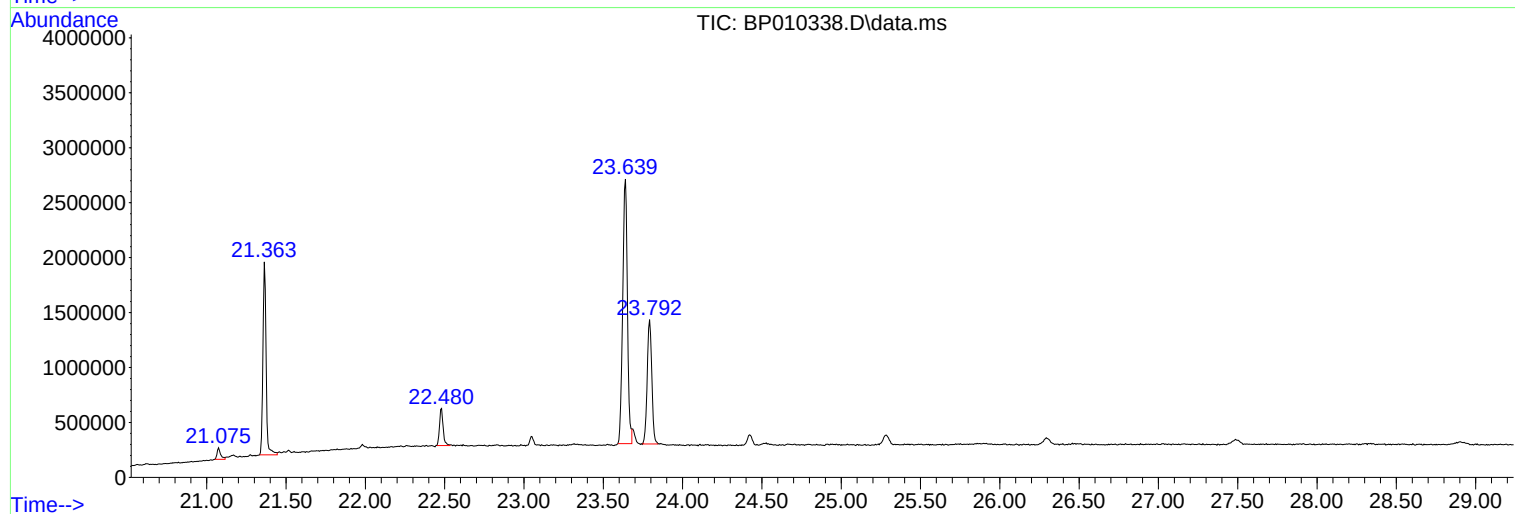
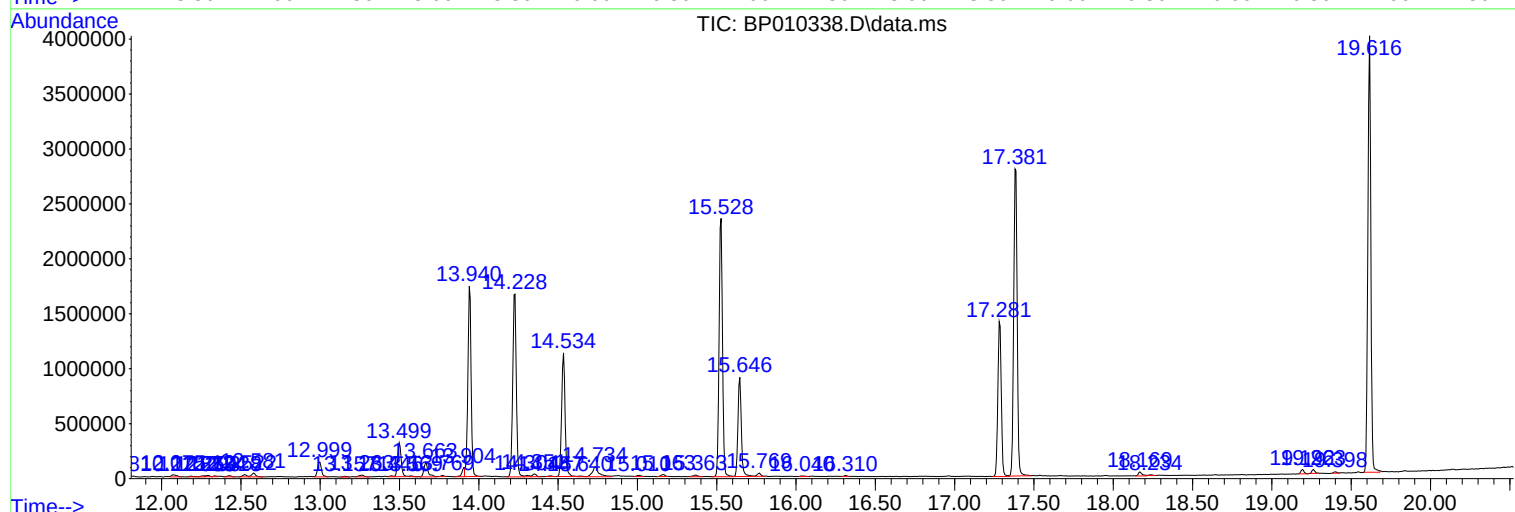
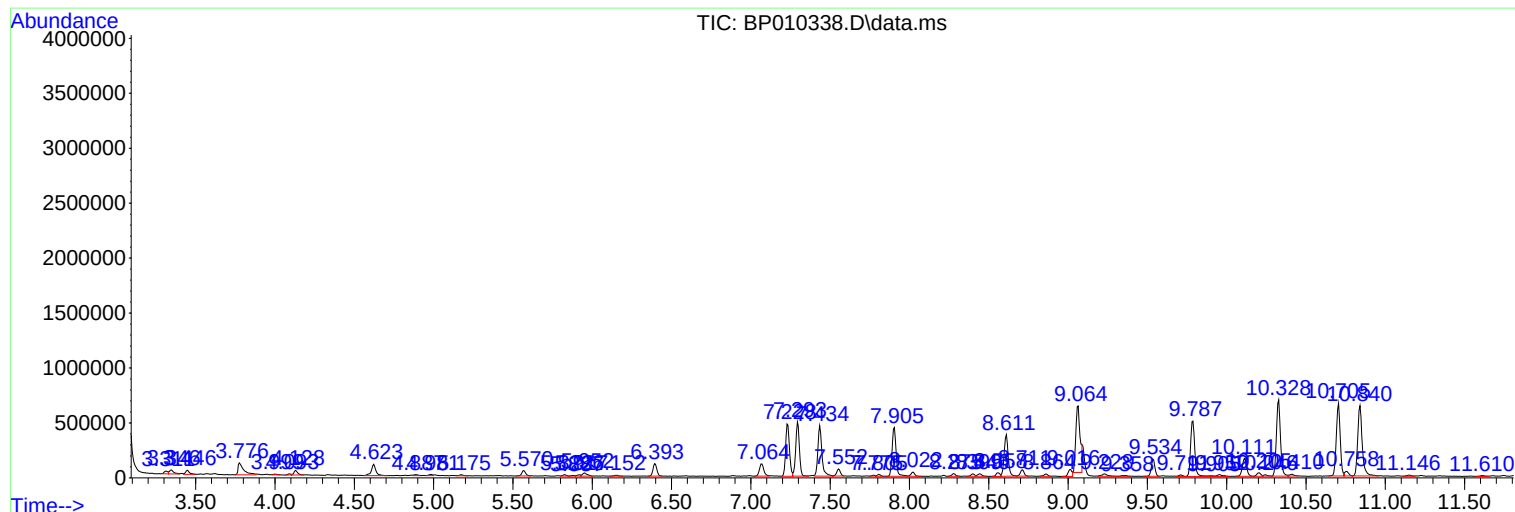
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION

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



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Instrument :
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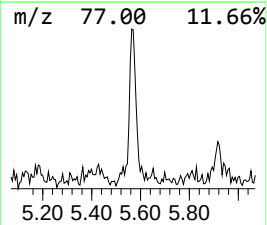
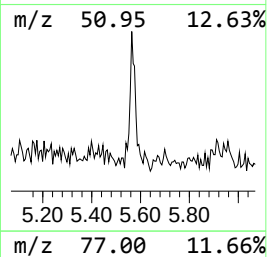
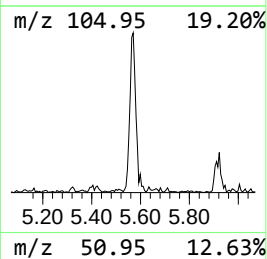
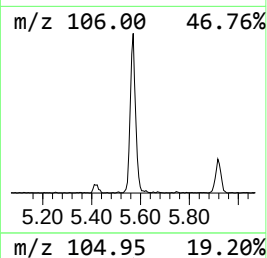
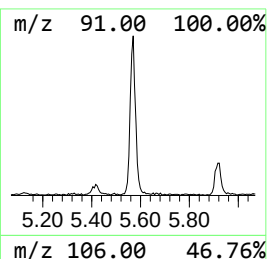
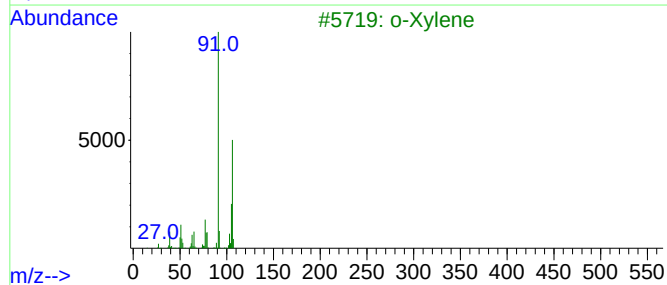
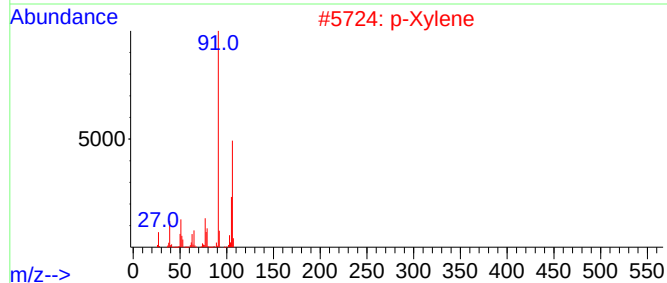
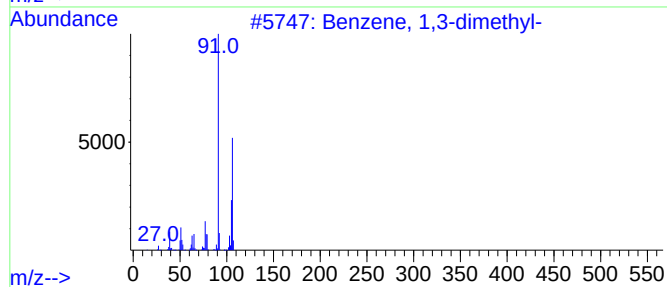
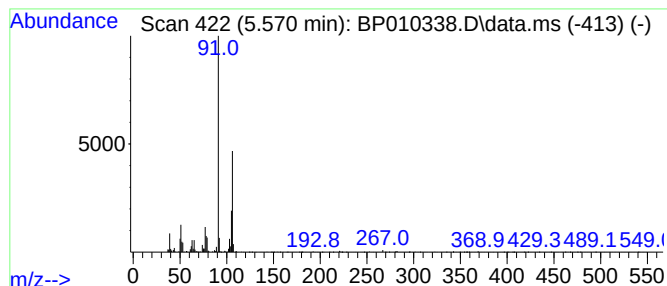
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,3-dimethyl- Concentration Rank 12

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.570	2.29 ng/ul	87660	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
2			p-Xylene	106	C8H10	000106-42-3	97
3			o-Xylene	106	C8H10	000095-47-6	97
4			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	97
5			Benzene, 1,3-dimethyl-	106	C8H10	000108-38-3	95



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ALS Vial : 62 Sample Multiplier: 1

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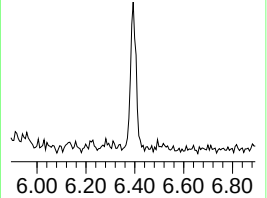
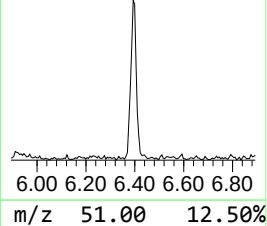
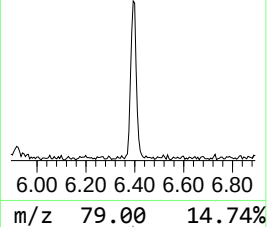
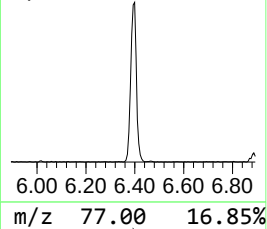
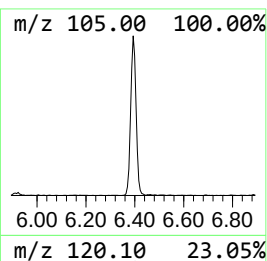
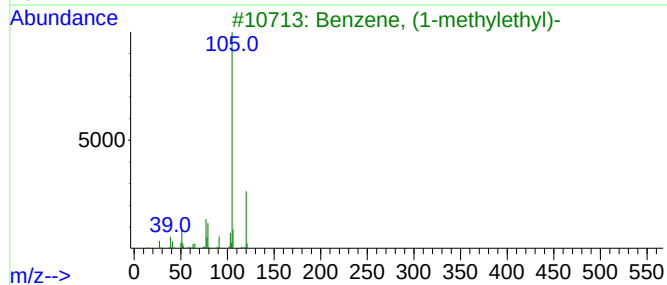
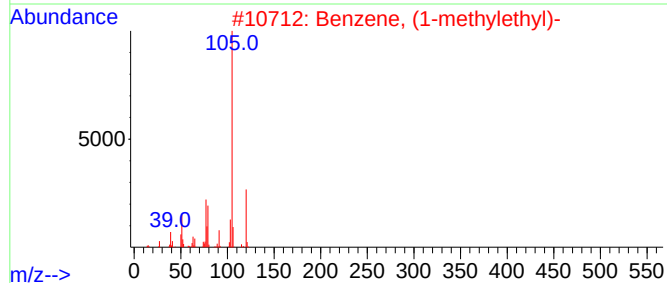
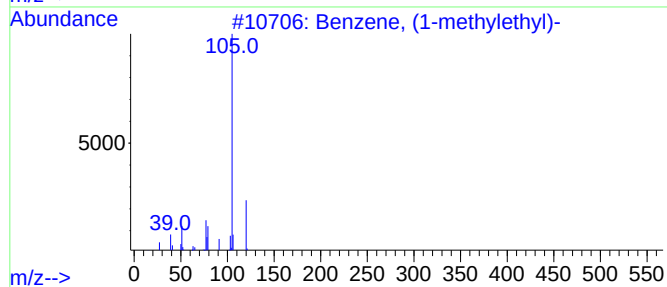
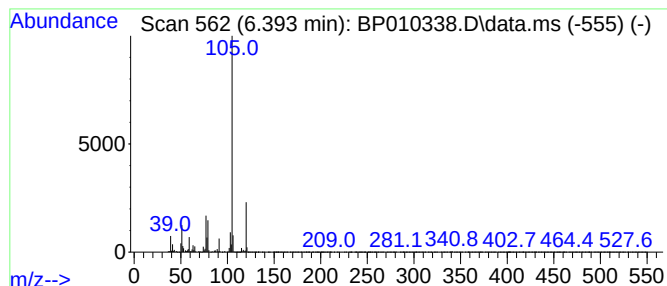
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Benzene, (1-methylethyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.393	5.05 ng/ul	193307	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
2			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	94
3			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	93
4			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	93
5			Benzene, (1-methylethyl)-	120	C9H12	000098-82-8	93



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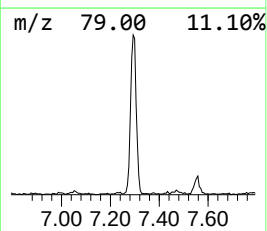
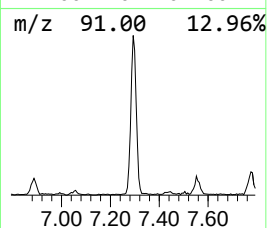
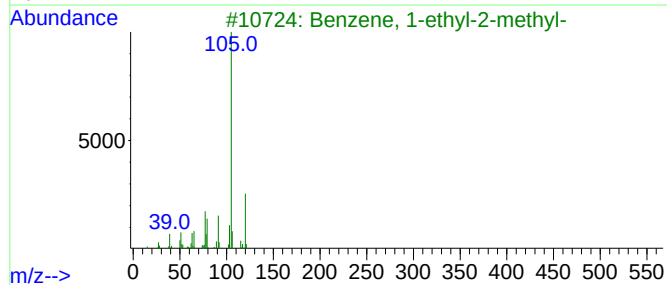
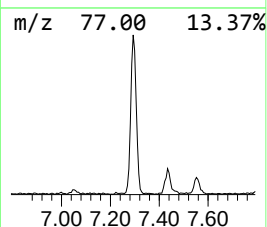
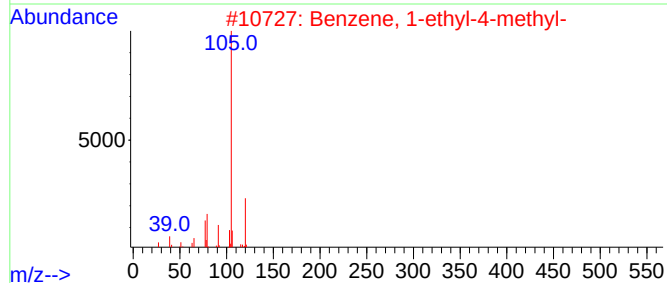
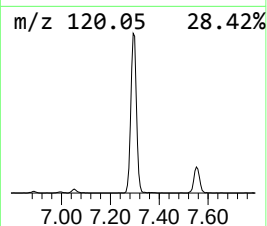
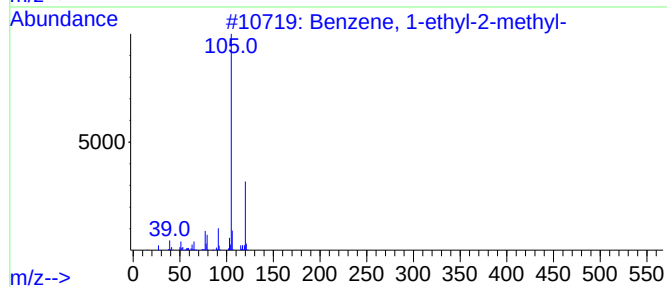
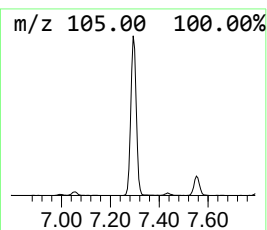
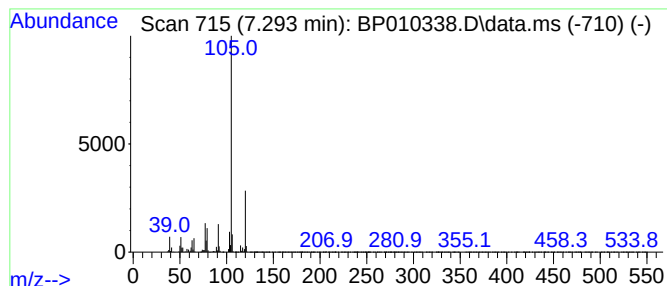
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1-ethyl-2-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.293	20.61 ng/ul	789194	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	91
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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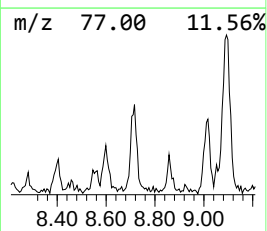
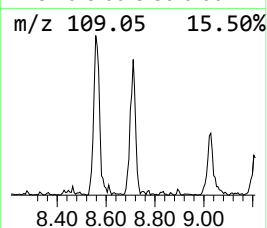
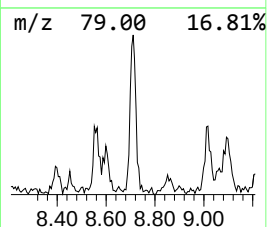
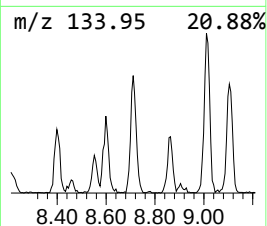
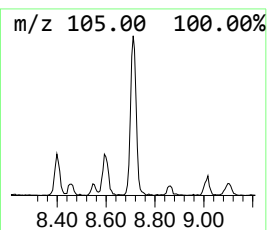
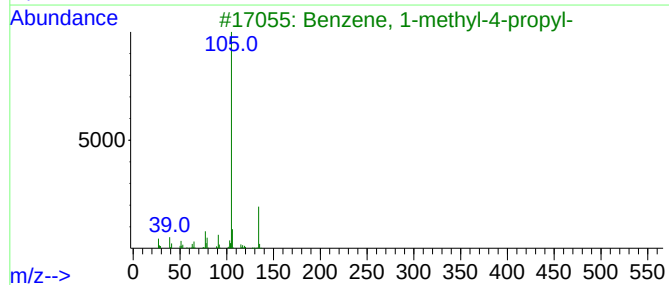
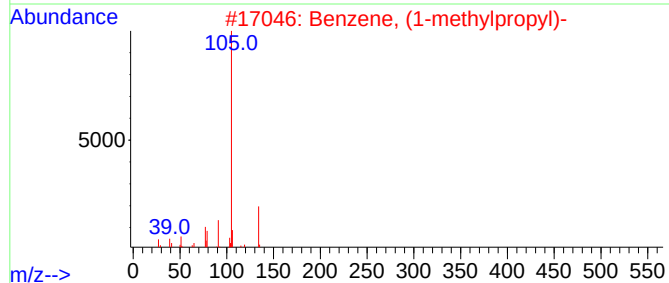
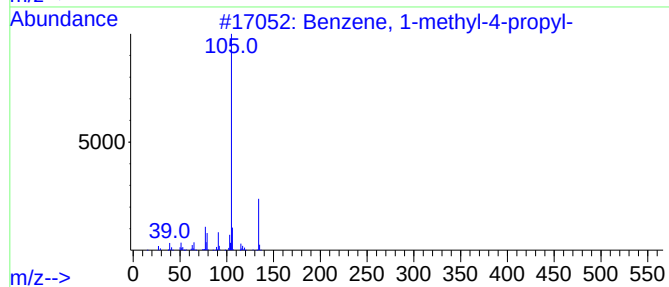
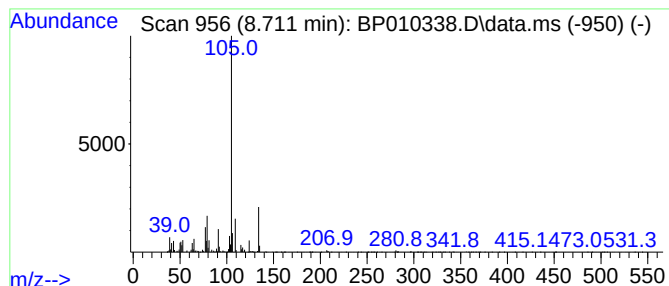
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1-methyl-4-propyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.711	2.51 ng/ul	95960	1,4-Dichlorobenzene-d4	7.905

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	93
2			Benzene, (1-methylpropyl)-	134	C10H14	000135-98-8	81
3			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	81
4			Benzene, 1-methyl-4-propyl-	134	C10H14	001074-55-1	76
5			Benzene, 1-methyl-2-propyl-	134	C10H14	001074-17-5	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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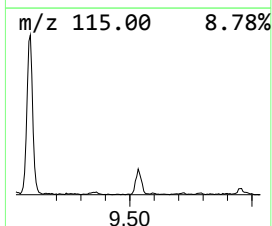
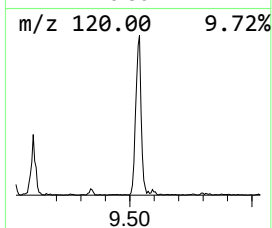
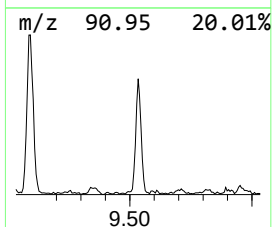
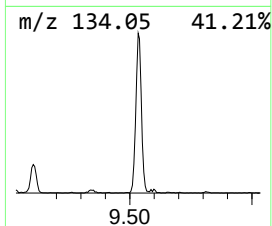
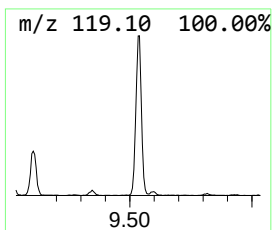
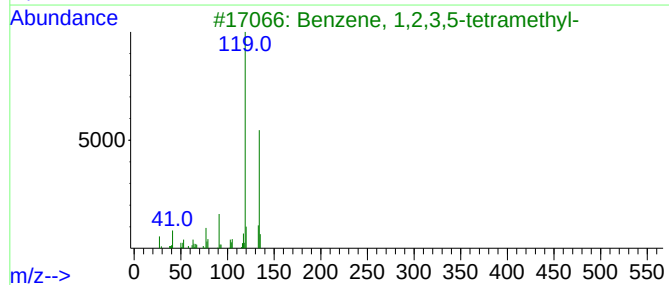
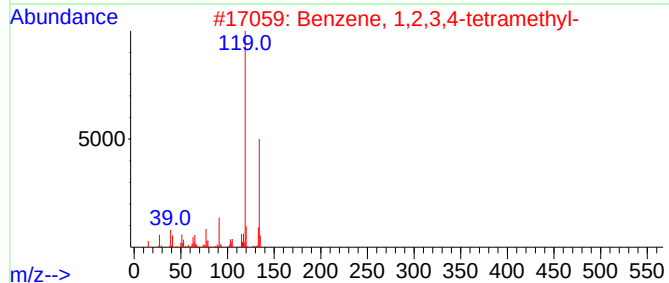
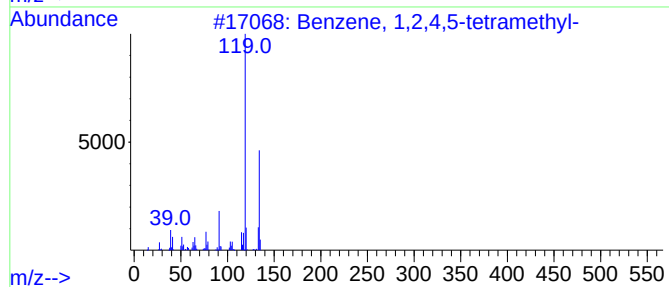
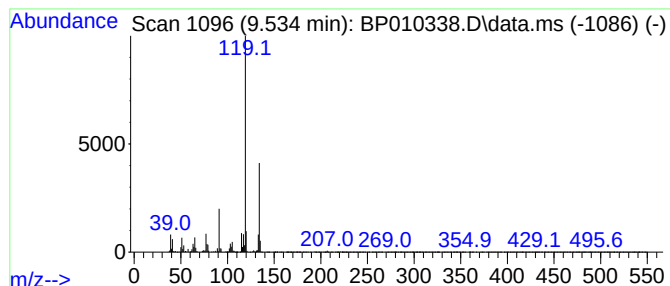
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.534	4.29 ng/ul	244481	Naphthalene-d8	10.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	97
2			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
4			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
5			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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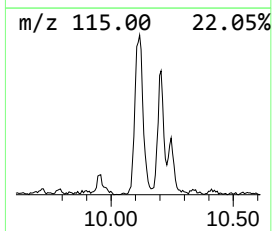
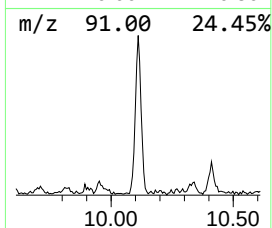
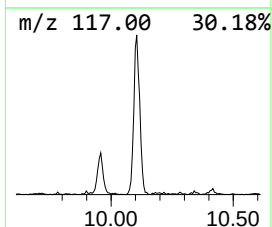
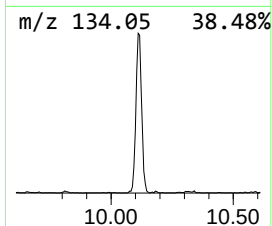
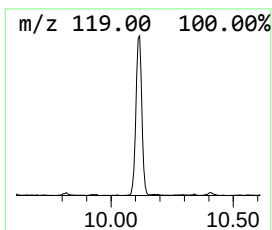
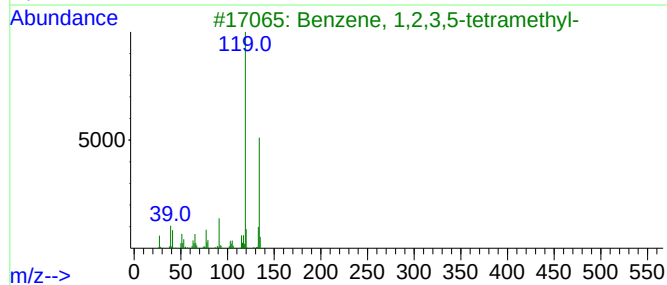
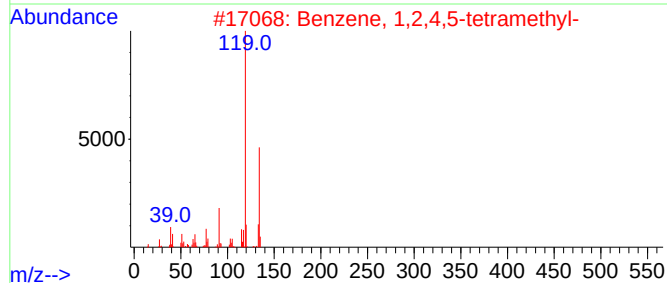
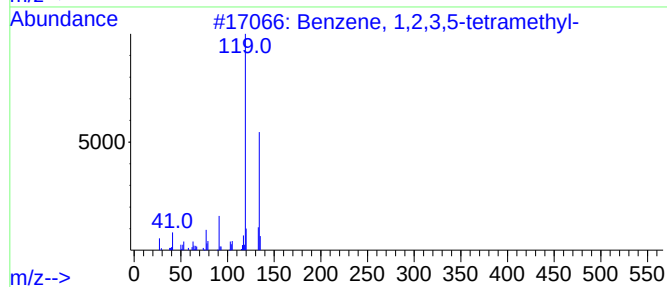
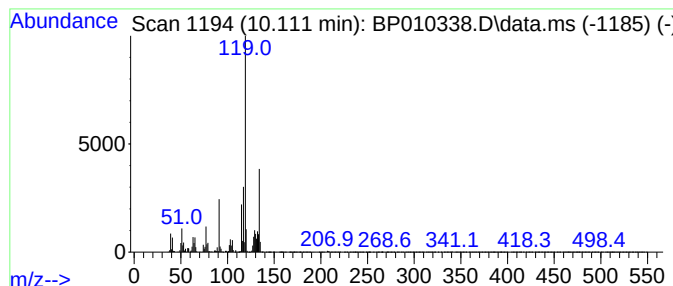
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.111	5.44 ng/ul	310070	Naphthalene-d8	10.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	70
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
3			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	70
4			p-Cymene	134	C10H14	000099-87-6	70
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
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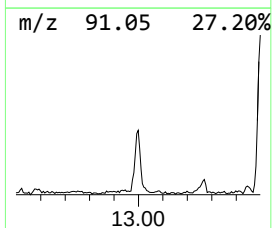
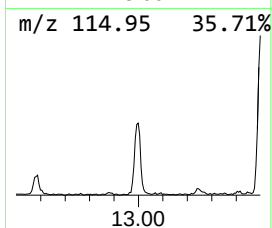
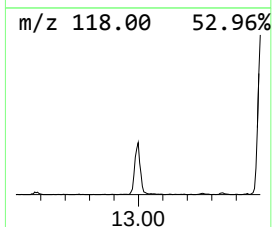
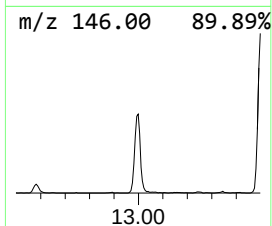
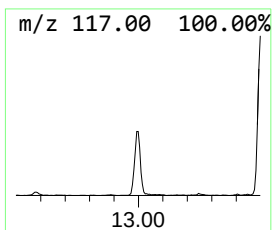
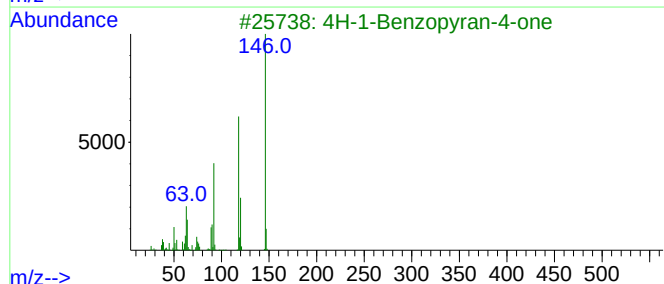
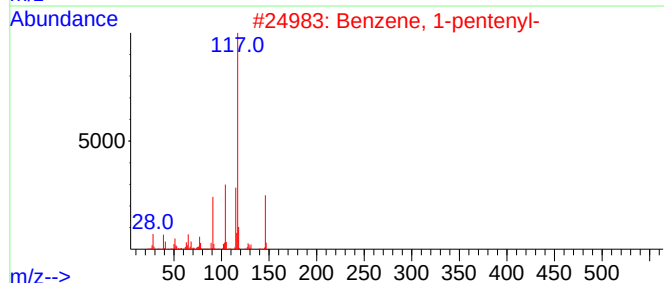
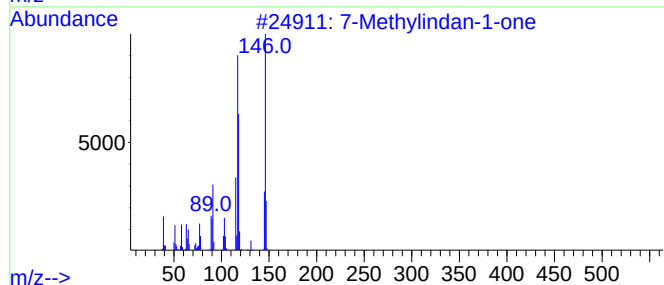
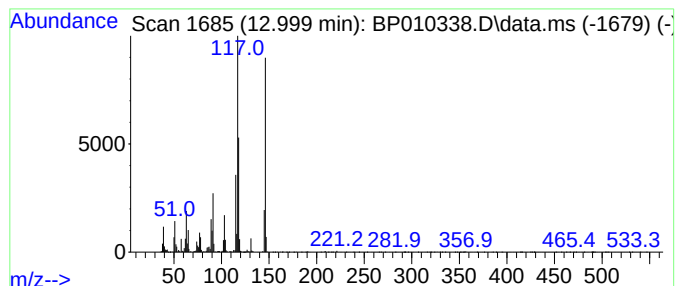
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 7-Methylindan-1-one Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.999	2.67 ng/ul	225700	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			7-Methylindan-1-one	146	C10H10O	039627-61-7	90
2			Benzene, 1-pentenyl-	146	C11H14	000826-18-6	58
3			4H-1-Benzopyran-4-one	146	C9H6O2	000491-38-3	50
4			Indane	118	C9H10	000496-11-7	50
5			Benzene, cyclopropyl-	118	C9H10	000873-49-4	50



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
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Sample : N2801-02
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Instrument :
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ClientSampleId :
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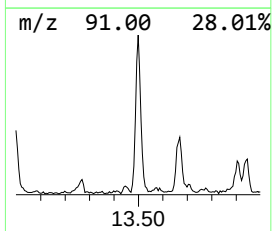
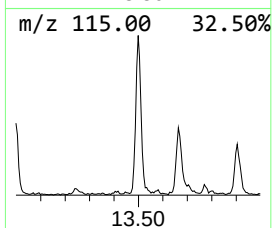
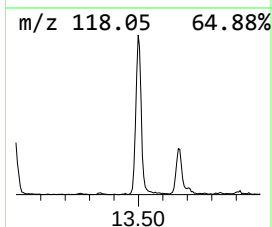
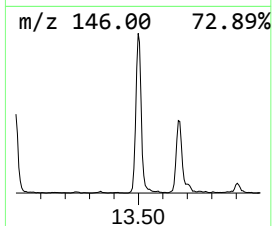
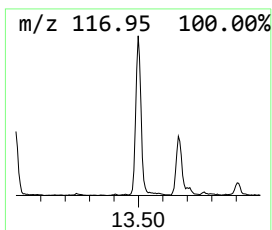
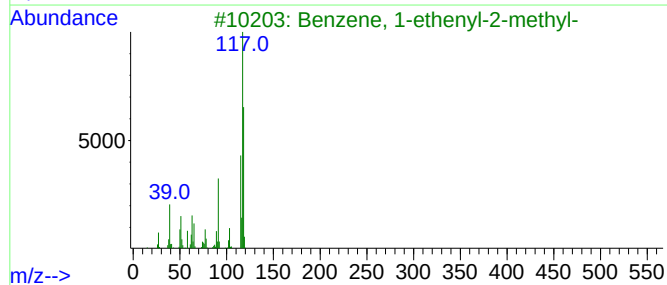
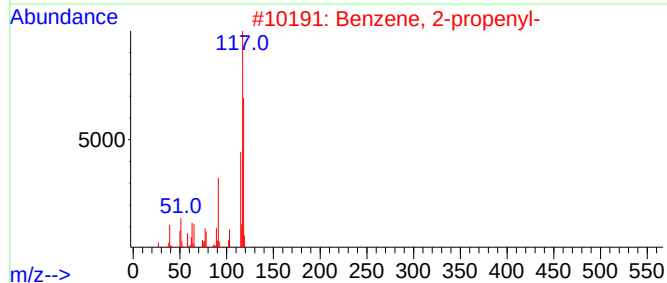
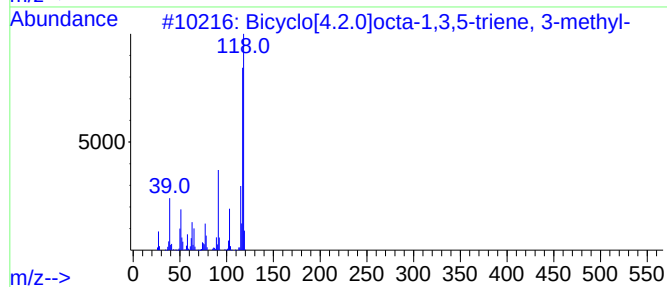
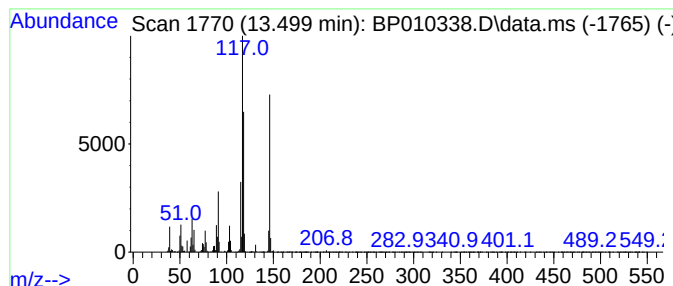
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.499	5.52 ng/ul	467079	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Bicyclo[4.2.0]octa-1,3,5-triene,...	118	C9H10	022250-74-4	86
2			Benzene, 2-propenyl-	118	C9H10	000300-57-2	83
3			Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	83
4			Benzene, 1-ethenyl-4-methyl-	118	C9H10	000622-97-9	83
5			(Z)-1-Phenylpropene	118	C9H10	000766-90-5	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
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ClientSampleId :
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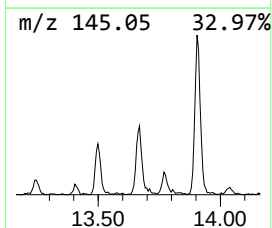
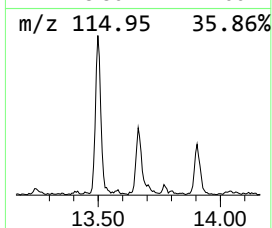
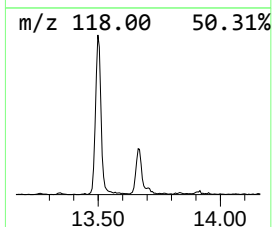
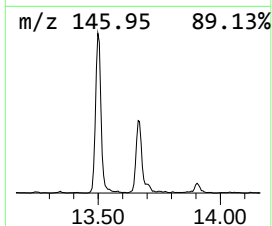
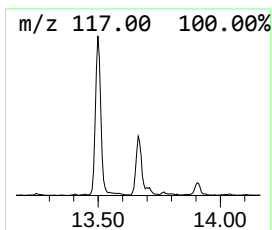
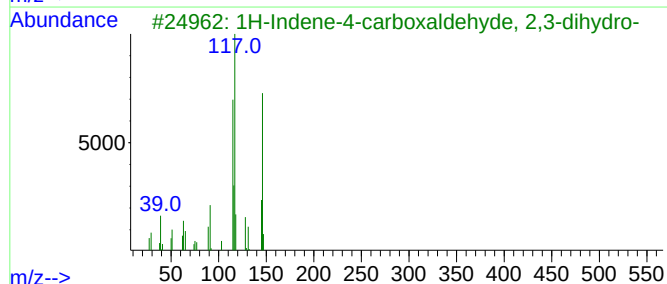
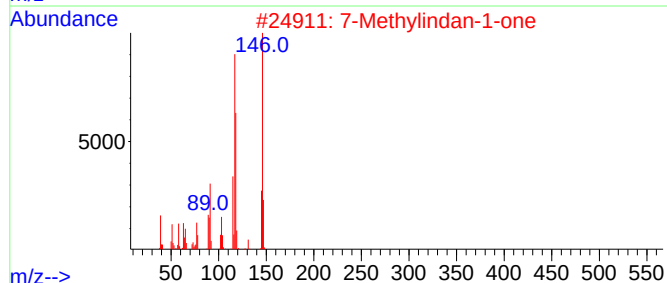
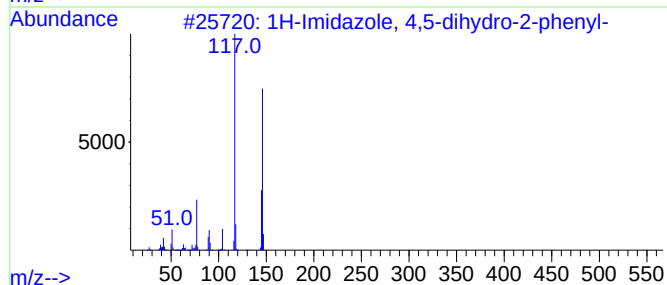
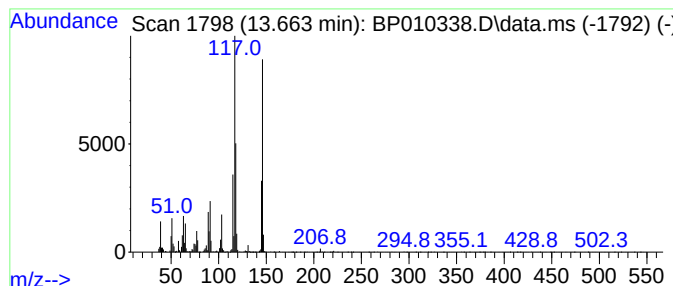
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1H-Imidazole, 4,5-dihydro-2... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.663	2.74 ng/ul	232257	Acenaphthene-d10	14.534

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Imidazole, 4,5-dihydro-2-phenyl-	146	C9H10N2	000936-49-2	76
2			7-Methylindan-1-one	146	C10H10O	039627-61-7	72
3			1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	64
4			1H-Indene-4-carboxaldehyde, 2,3-...	146	C10H10O	051932-70-8	64
5			Benzene, 1-pentenyl-	146	C11H14	000826-18-6	58



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
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Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

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ClientSampleId :
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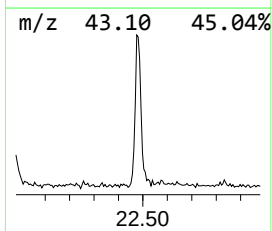
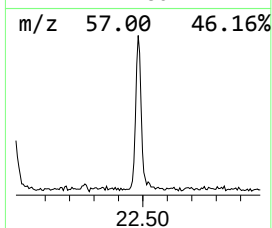
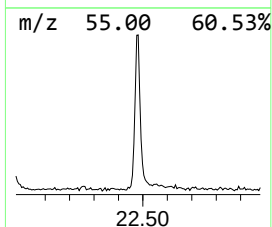
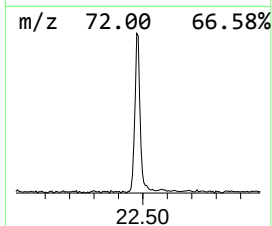
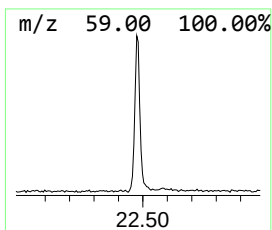
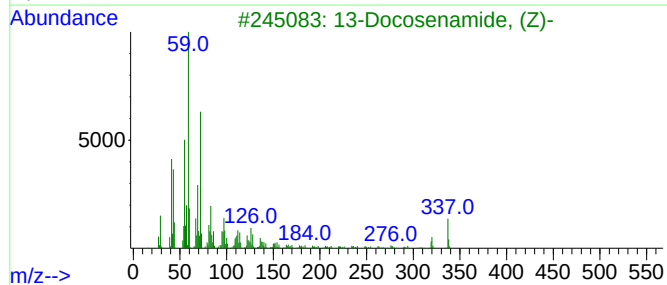
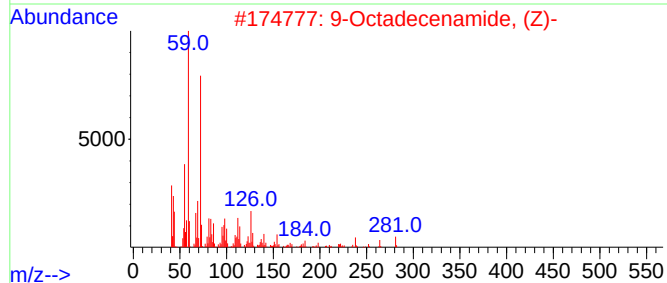
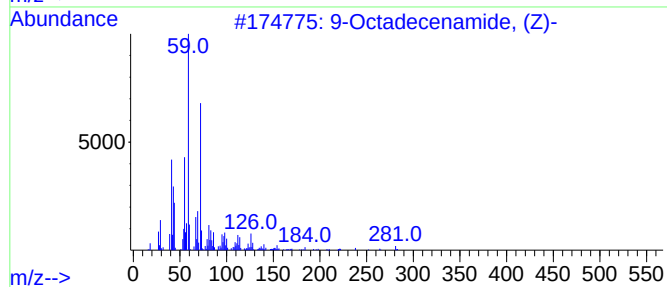
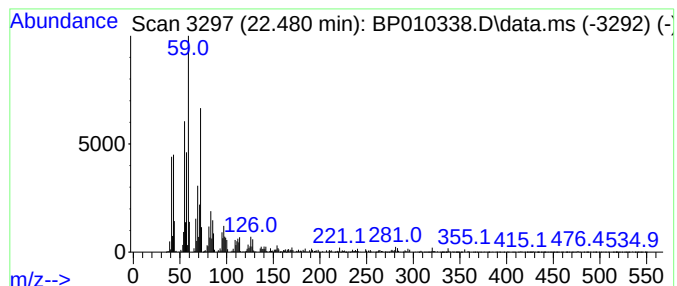
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 9-Octadecenamide, (Z)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
22.480	4.41 ng/ul	529521	Chrysene-d12	21.363

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	98
2		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	93
3		13-Docosenamide, (Z)-	337	C22H43NO	000112-84-5	86
4		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	83
5		9-Octadecenamide, (Z)-	281	C18H35NO	000301-02-0	78



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP051122\
Data File : BP010338.D
Acq On : 12 May 2022 23:38
Operator : CG/JU
Sample : N2801-02
Misc :
ALS Vial : 62 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
EXLB5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP051022.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,3-di...	5.570	2.3	ng/ul	87660	1	7.905	766010	20.0
Benzene, (1-met...	6.393	5.0	ng/ul	193307	1	7.905	766010	20.0
Benzene, 1-ethy...	7.293	20.6	ng/ul	789194	1	7.905	766010	20.0
Benzene, 1-meth...	8.711	2.5	ng/ul	95960	1	7.905	766010	20.0
Benzene, 1,2,4,...	9.534	4.3	ng/ul	244481	2	10.705	1140560	20.0
Benzene, 1,2,3,...	10.111	5.4	ng/ul	310070	2	10.705	1140560	20.0
7-Methylindan-1...	12.999	2.7	ng/ul	225700	3	14.534	1693700	20.0
Bicyclo[4.2.0]o...	13.499	5.5	ng/ul	467079	3	14.534	1693700	20.0
1H-Imidazole, 4...	13.663	2.7	ng/ul	232257	3	14.534	1693700	20.0
9-Octadecenamid...	22.480	4.4	ng/ul	529521	5	21.363	2403560	20.0