

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
 Data File : BP014438.D
 Acq On : 31 Mar 2023 12:00
 Operator : CG/JU
 Sample : 02131-02DL 5X
 Misc :
 ALS Vial : 47 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0QG7DL

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-BP032823.M
 Title : SVOA CALIBRATION

Signal : TIC: BP014438.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.864	113	115	126	rBV	17975	47344	0.22%	0.083%
2	4.146	159	163	168	rVB	22412	31858	0.15%	0.056%
3	5.622	409	414	423	rBV	32297	68546	0.31%	0.121%
4	5.993	471	477	482	rBV3	37382	67244	0.31%	0.118%
5	6.046	482	486	496	rVB	27015	42880	0.20%	0.075%
6	7.087	657	663	668	rBV	38001	62018	0.28%	0.109%
7	7.146	668	673	676	rVV	24254	41026	0.19%	0.072%
8	7.187	676	680	681	rVV	23963	32669	0.15%	0.057%
9	7.222	682	686	695	rVB	55253	92856	0.43%	0.163%
10	7.340	700	706	711	rBV	144839	232527	1.07%	0.409%
11	7.387	711	714	719	rVB2	27223	43838	0.20%	0.077%
12	7.546	734	741	748	rBV	121709	215380	0.99%	0.379%
13	7.610	748	752	754	rVV	50488	69479	0.32%	0.122%
14	7.652	754	759	766	rVB	205804	327597	1.50%	0.576%
15	7.728	766	772	779	rBV	18224	34190	0.16%	0.060%
16	8.010	815	820	828	rVV	639630	1036362	4.75%	1.823%
17	8.122	834	839	846	rVB	125652	208583	0.96%	0.367%
18	8.387	877	884	889	rBV	75794	117787	0.54%	0.207%
19	8.552	906	912	921	rVB2	169294	297261	1.36%	0.523%
20	8.646	921	928	934	rBV2	77358	137454	0.63%	0.242%
21	8.734	934	943	952	rVV	85506	178578	0.82%	0.314%
22	8.816	953	957	963	rVV	17110	36842	0.17%	0.065%
23	8.881	963	968	975	rVB5	16847	36820	0.17%	0.065%
24	8.963	975	982	986	rBV	40626	64269	0.29%	0.113%
25	9.016	986	991	997	rVB	43351	64434	0.30%	0.113%
26	9.116	997	1008	1014	rBV2	83405	160372	0.74%	0.282%
27	9.193	1014	1021	1038	rVB3	190101	455292	2.09%	0.801%
28	9.369	1046	1051	1059	rVB7	17518	38838	0.18%	0.068%
29	9.457	1059	1066	1070	rBV2	32322	59123	0.27%	0.104%
30	9.510	1070	1075	1080	rVB8	13561	29515	0.14%	0.052%
31	9.646	1093	1098	1103	rBV	60672	91721	0.42%	0.161%
32	9.704	1103	1108	1116	rVV	82306	132867	0.61%	0.234%
33	9.916	1138	1144	1156	rVB	132838	263939	1.21%	0.464%
34	10.016	1156	1161	1166	rVV3	21029	33032	0.15%	0.058%
35	10.075	1166	1171	1181	rVB	48377	85983	0.39%	0.151%
36	10.228	1191	1197	1205	rBV4	110895	217802	1.00%	0.383%

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Integrator: RTE

Smoothing : OFF

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
 Title : SVOA CALIBRATION

37	10.340	1211	1216	1225	rBV5	19987	46186	0.21%	0.081%
38	10.469	1231	1238	1245	rBV	181984	364483	1.67%	0.641%
39	10.534	1245	1249	1258	rVB5	40575	83100	0.38%	0.146%
40	10.787	1287	1292	1294	rBV	32964	49114	0.23%	0.086%

41	10.834	1294	1300	1304	rVV	969400	1691016	7.75%	2.975%
42	10.887	1304	1309	1317	rVV	2621070	4362882	20.01%	7.674%
43	10.993	1321	1327	1341	rVB2	99681	279892	1.28%	0.492%
44	11.946	1483	1489	1497	rBV6	14641	38919	0.18%	0.068%
45	12.234	1530	1538	1544	rBV3	28416	60417	0.28%	0.106%

46	12.387	1559	1564	1572	rBV5	12035	31853	0.15%	0.056%
47	12.493	1576	1582	1595	rVB	475862	762900	3.50%	1.342%
48	12.710	1607	1619	1629	rBV	436867	704870	3.23%	1.240%
49	13.469	1743	1748	1749	rBV	33680	38093	0.17%	0.067%
50	13.498	1749	1753	1760	rVB2	116738	181292	0.83%	0.319%

51	13.693	1773	1786	1789	rBV3	65149	147217	0.68%	0.259%
52	13.828	1802	1809	1810	rBV2	90408	141794	0.65%	0.249%
53	13.987	1829	1836	1841	rBV	196790	362371	1.66%	0.637%
54	14.040	1841	1845	1847	rVV	137894	203787	0.93%	0.358%
55	14.069	1847	1850	1857	rVV	535536	754558	3.46%	1.327%

56	14.222	1871	1876	1879	rBV	104789	155765	0.71%	0.274%
57	14.251	1879	1881	1886	rVB2	47822	58279	0.27%	0.103%
58	14.357	1893	1899	1915	rBV2	522815	943487	4.33%	1.660%
59	14.540	1926	1930	1939	rBV2	52950	72045	0.33%	0.127%
60	14.663	1939	1951	1958	rBV	1591785	2547913	11.68%	4.482%

61	14.740	1959	1964	1970	rVV3	55331	124096	0.57%	0.218%
62	14.875	1981	1987	1994	rBV2	38999	93960	0.43%	0.165%
63	14.957	1997	2001	2008	rBV4	16624	35383	0.16%	0.062%
64	15.063	2013	2019	2026	rVV3	89322	169746	0.78%	0.299%
65	15.134	2026	2031	2040	rVV2	55890	97288	0.45%	0.171%

66	15.216	2040	2045	2050	rVB3	74870	120353	0.55%	0.212%
67	15.298	2050	2059	2070	rBV2	767061	1356626	6.22%	2.386%
68	15.445	2077	2084	2088	rBV3	43659	101743	0.47%	0.179%
69	15.492	2088	2092	2098	rVB2	97053	132233	0.61%	0.233%
70	15.598	2105	2110	2111	rBV4	21536	30766	0.14%	0.054%

71	15.663	2111	2121	2126	rVV	692346	1073454	4.92%	1.888%
72	15.710	2126	2129	2133	rVB3	58465	94477	0.43%	0.166%
73	15.792	2138	2143	2149	rBV2	168263	304578	1.40%	0.536%
74	15.875	2154	2157	2163	rVB4	27123	45960	0.21%	0.081%
75	16.016	2173	2181	2190	rVB6	35981	105756	0.48%	0.186%

76	16.157	2202	2205	2213	rVB5	24074	42367	0.19%	0.075%
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 Title : SVOA CALIBRATION

77	16.357	2235	2239	2242	rBV	74243	112632	0.52%	0.198%
78	16.704	2294	2298	2305	rBV6	42106	99187	0.45%	0.174%
79	16.775	2306	2310	2318	rVV2	46800	87751	0.40%	0.154%
80	16.875	2324	2327	2331	rVB4	31446	40403	0.19%	0.071%
81	16.934	2332	2337	2340	rBV5	21523	47382	0.22%	0.083%
82	16.963	2340	2342	2352	rVB10	23844	51968	0.24%	0.091%
83	17.081	2355	2362	2372	rBV	14424452	21806027	100.00%	38.357%
84	17.157	2373	2375	2380	rVV3	103068	167663	0.77%	0.295%
85	17.210	2381	2384	2387	rVV4	62500	95775	0.44%	0.168%
86	17.251	2388	2391	2396	rVB	48572	69186	0.32%	0.122%
87	17.428	2415	2421	2426	rBV	1958782	2810564	12.89%	4.944%
88	17.469	2426	2428	2433	rVV	169352	219824	1.01%	0.387%
89	17.528	2433	2438	2449	rVB2	757212	1114719	5.11%	1.961%
90	17.898	2497	2501	2505	rVB	34527	43084	0.20%	0.076%
91	18.292	2564	2568	2572	rBV	86687	127500	0.58%	0.224%
92	18.339	2573	2576	2582	rVB	63010	85337	0.39%	0.150%
93	18.469	2594	2598	2603	rBV3	65254	113084	0.52%	0.199%
94	18.516	2603	2606	2610	rVB	49397	59592	0.27%	0.105%
95	19.175	2714	2718	2721	rBV2	71510	98846	0.45%	0.174%
96	19.657	2796	2800	2804	rBV2	49739	62156	0.29%	0.109%
97	19.757	2812	2817	2829	rVB	875716	1207090	5.54%	2.123%
98	21.516	3111	3116	3126	rBV	1621062	2201957	10.10%	3.873%
99	23.868	3511	3516	3525	rVB3	479075	988796	4.53%	1.739%
100	24.033	3538	3544	3555	rVB2	936288	1972147	9.04%	3.469%

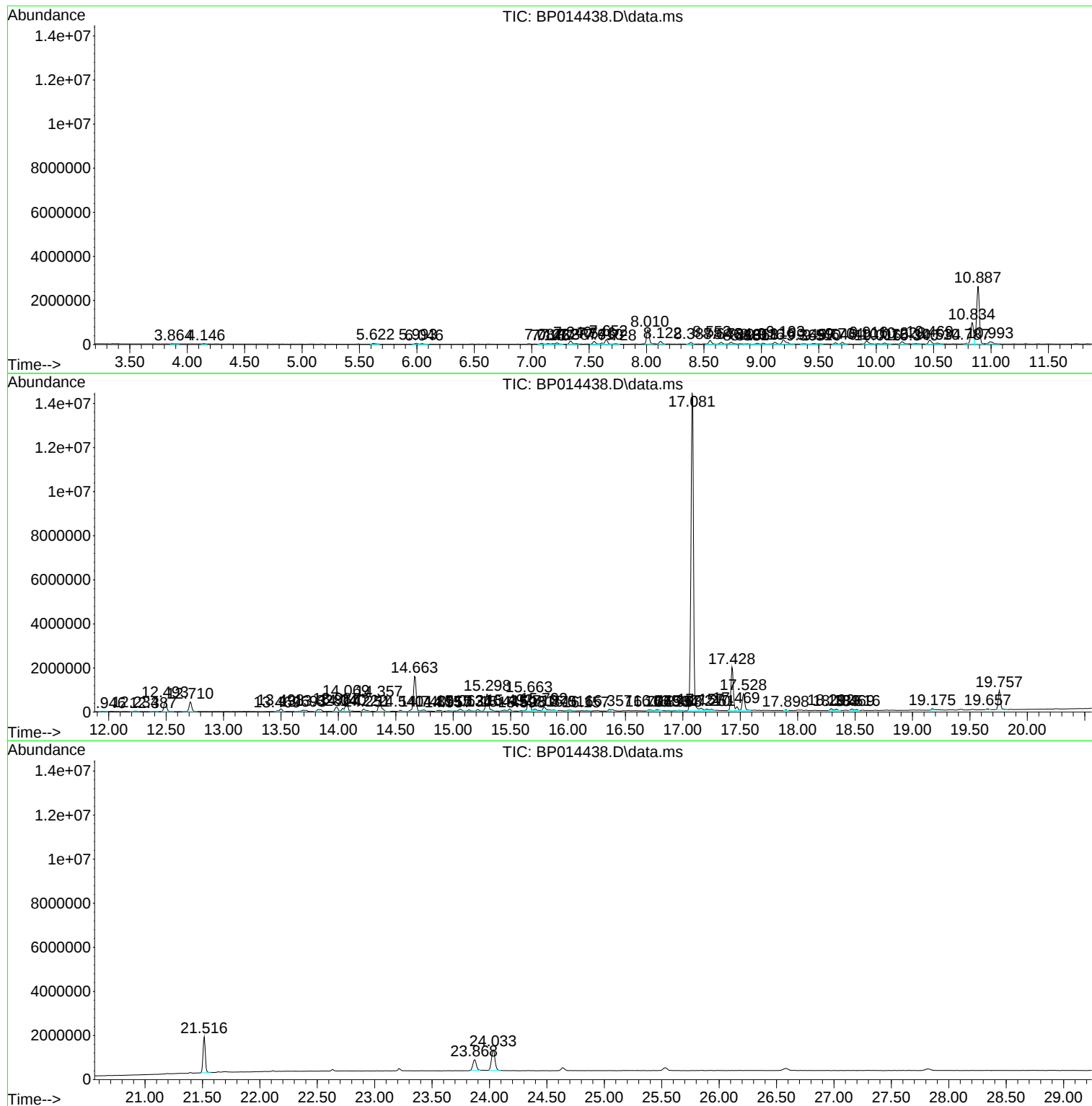
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Instrument :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
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ALS Vial : 47 Sample Multiplier: 1

Instrument :
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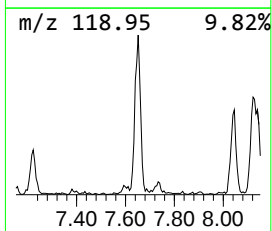
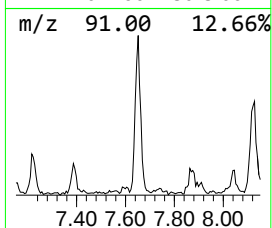
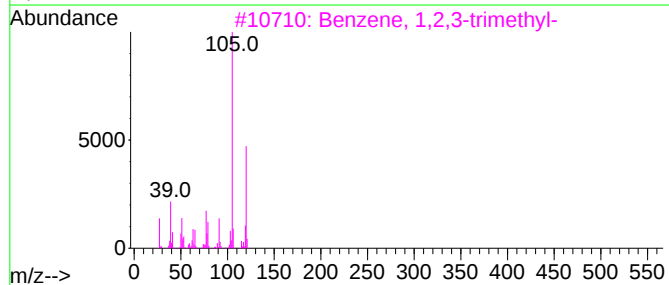
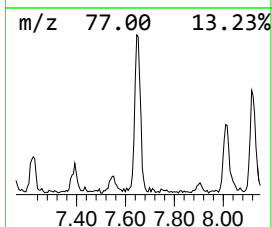
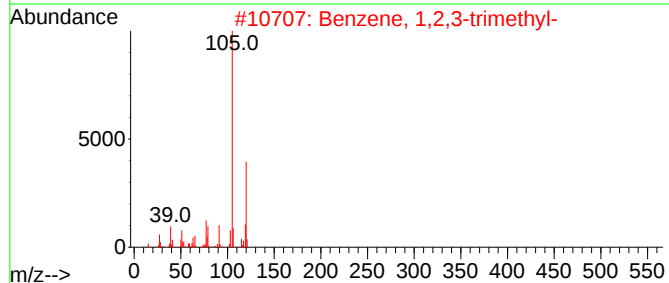
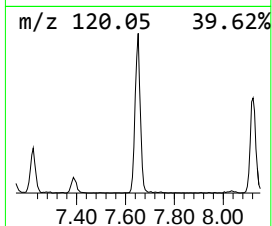
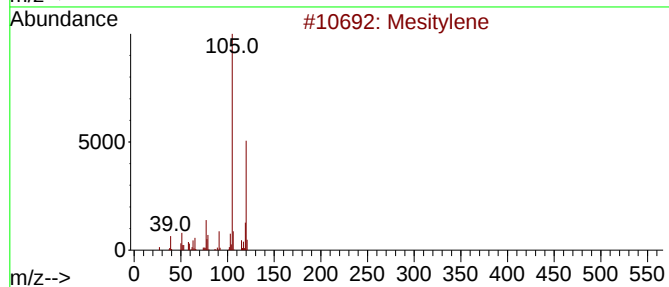
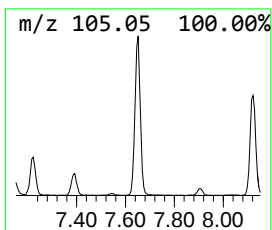
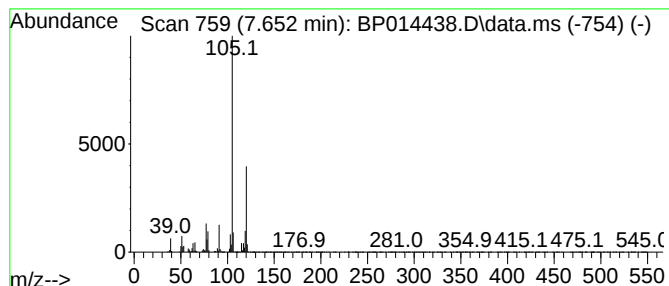
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Mesitylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.652	6.32 ng/ul	327597	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Mesitylene	120	C9H12	000108-67-8	97
2			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	95
3			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
4			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
5			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91



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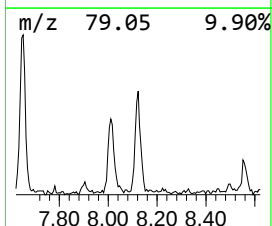
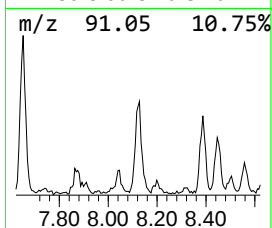
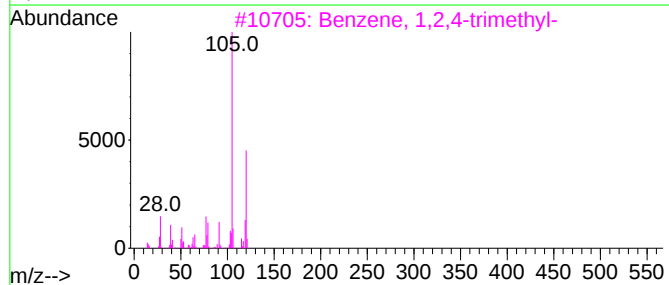
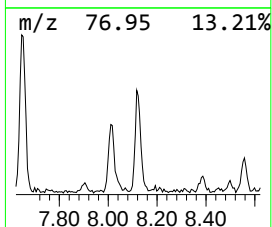
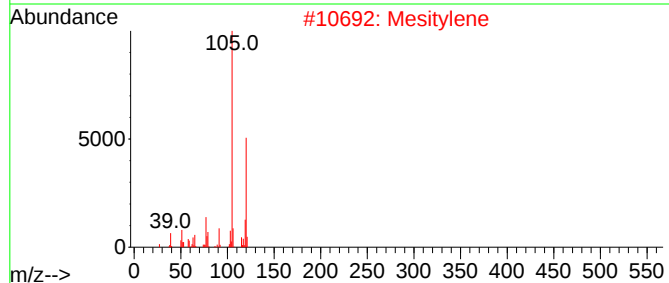
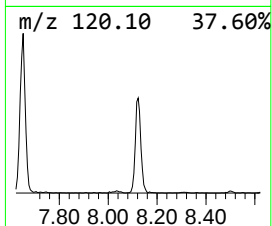
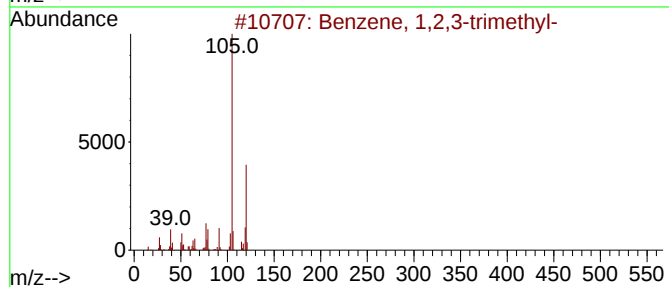
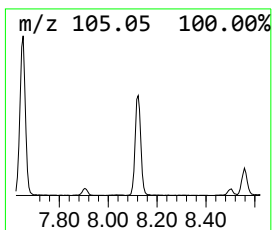
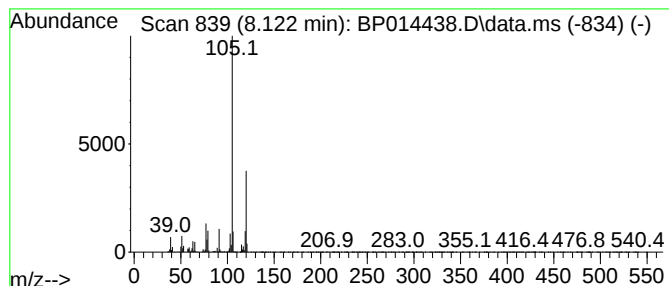
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.122	4.03 ng/ul	208583	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Mesitylene	120	C9H12	000108-67-8	95
3			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
4			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	94
5			Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	93



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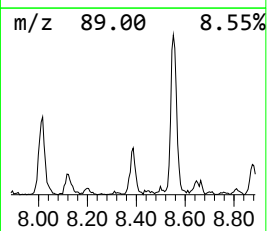
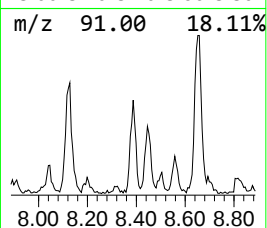
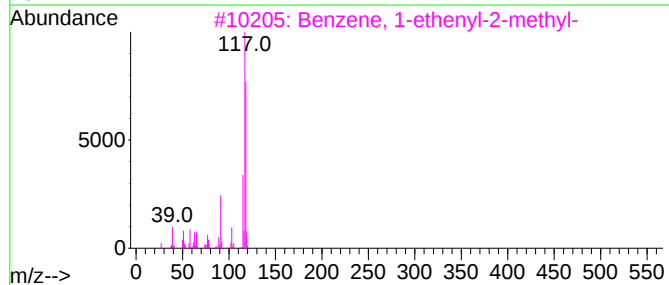
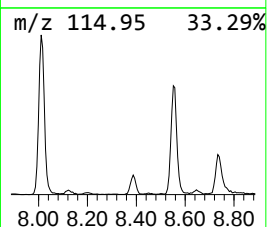
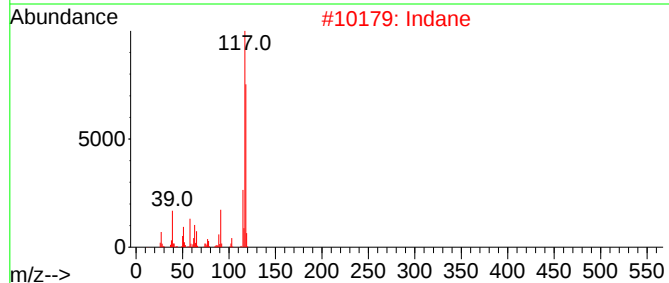
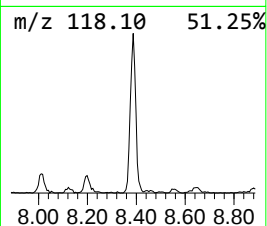
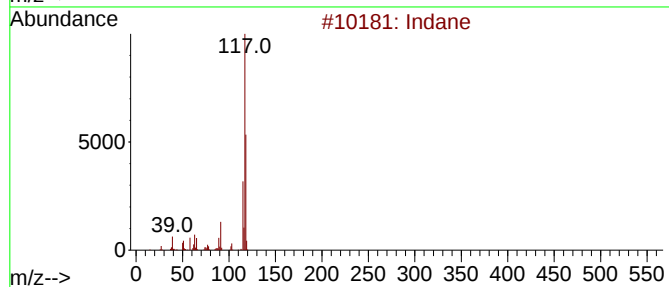
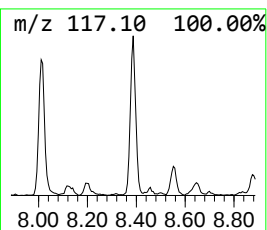
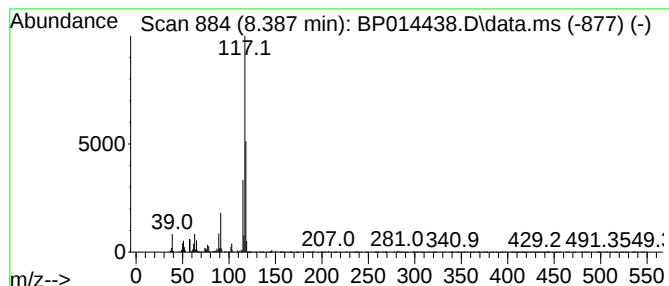
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 Indane Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.387	2.27 ng/ul	117787	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	94
2	Indane			118	C9H10	000496-11-7	87
3	Benzene, 1-ethenyl-2-methyl-			118	C9H10	000611-15-4	74
4	Benzene, cyclopropyl-			118	C9H10	000873-49-4	74
5	(Z)-1-Phenylpropene			118	C9H10	000766-90-5	72



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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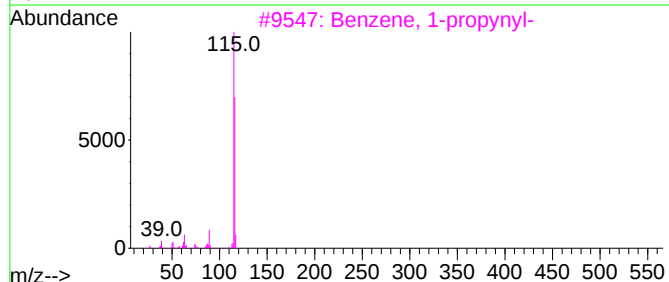
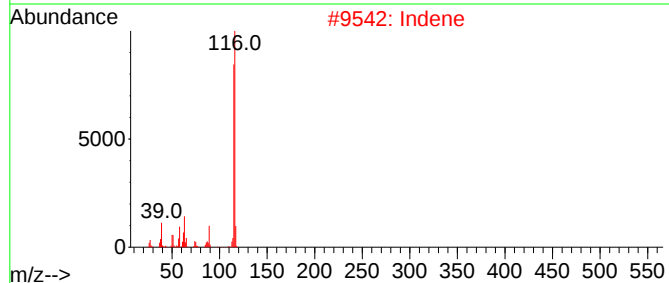
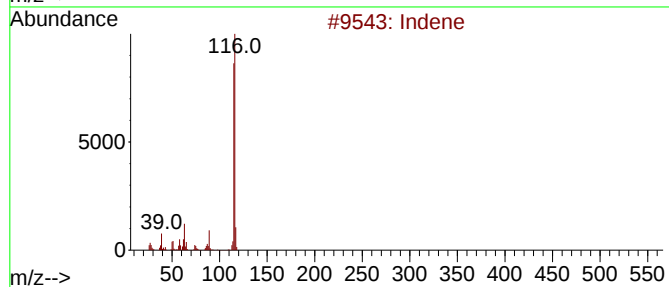
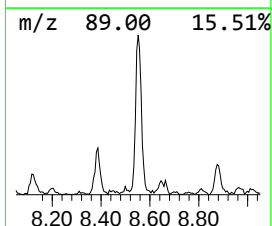
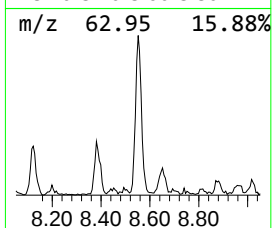
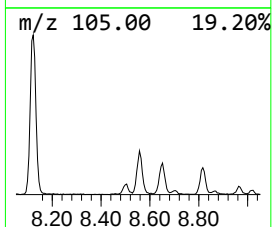
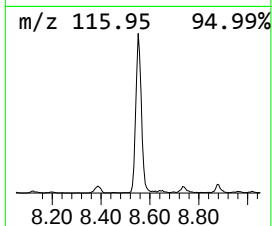
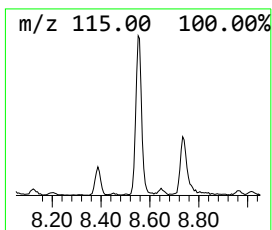
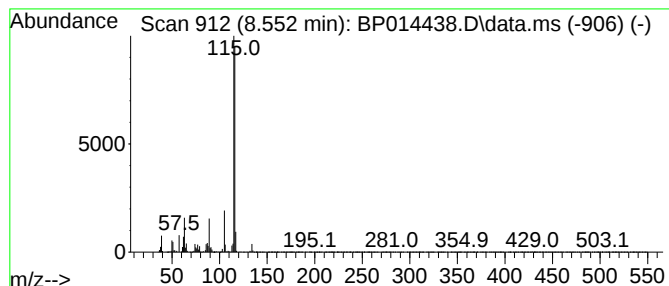
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Indene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.552	5.74 ng/ul	297261	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indene			116	C9H8	000095-13-6	94
2	Indene			116	C9H8	000095-13-6	94
3	Benzene, 1-propynyl-			116	C9H8	000673-32-5	87
4	Benzene, 1-ethynyl-4-methyl-			116	C9H8	000766-97-2	87
5	Benzene, 1-propynyl-			116	C9H8	000673-32-5	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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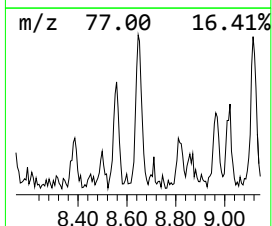
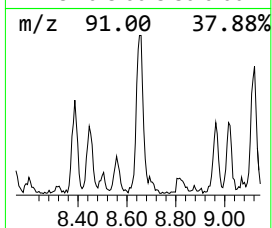
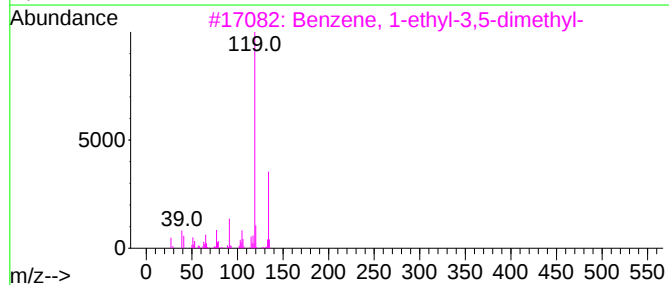
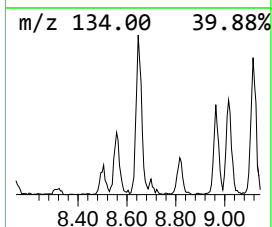
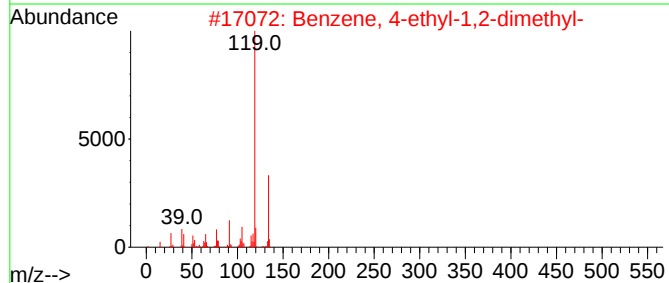
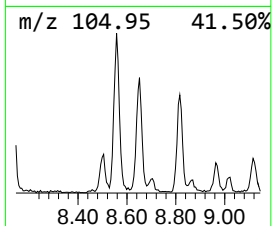
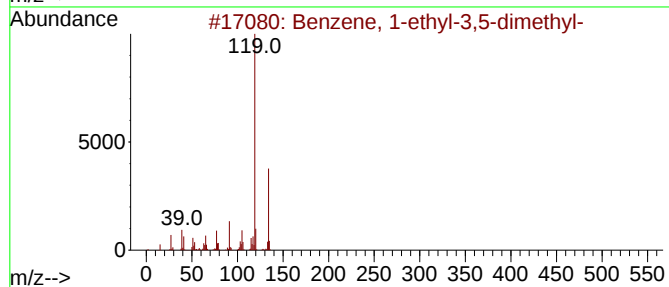
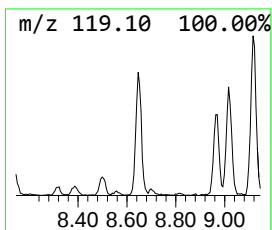
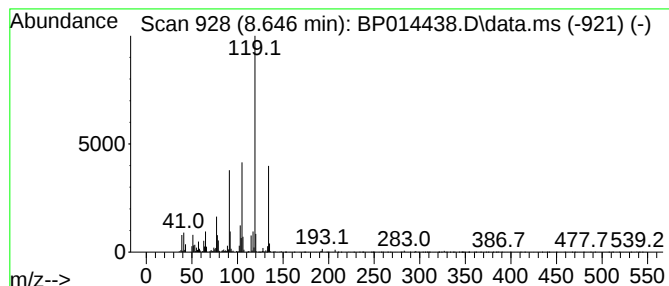
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 Benzene, 1-ethyl-3,5-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.646	2.65 ng/ul	137454	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	94
2			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
3			Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	91
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	91
5			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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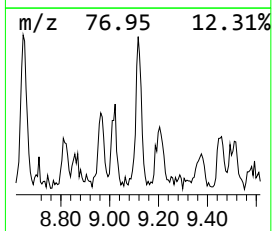
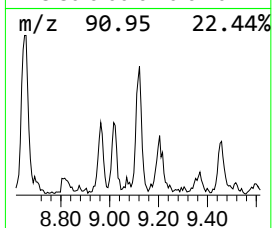
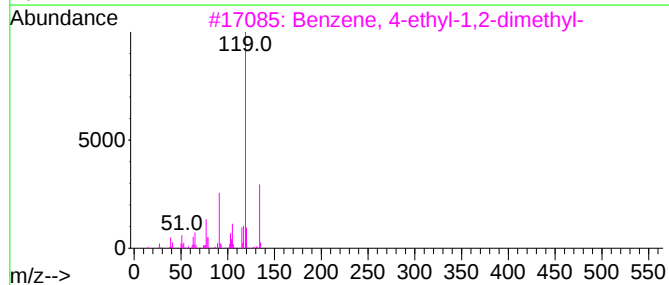
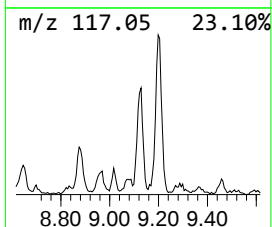
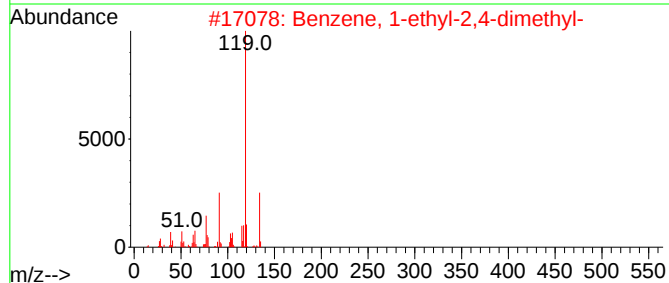
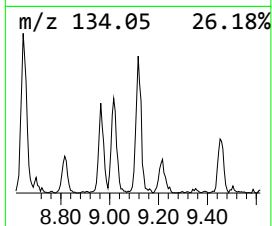
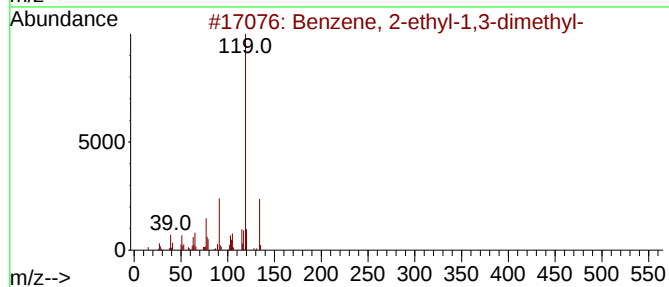
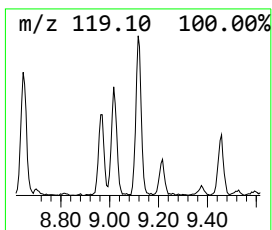
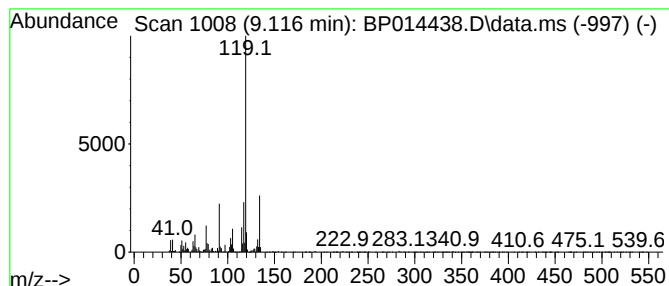
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.116	3.09 ng/ul	160372	1,4-Dichlorobenzene-d4	8.010

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
2			Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	93
5			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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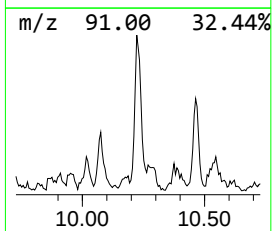
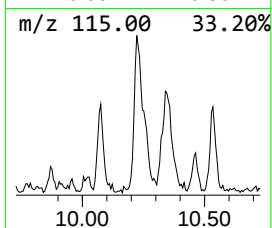
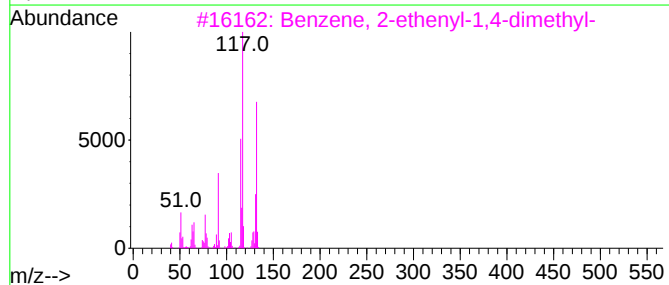
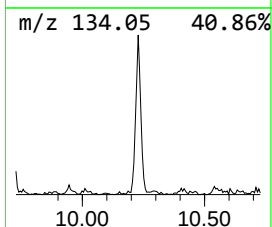
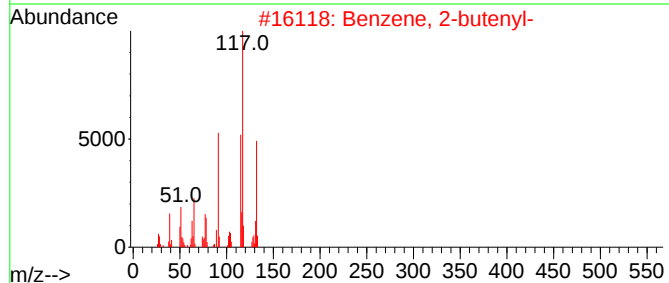
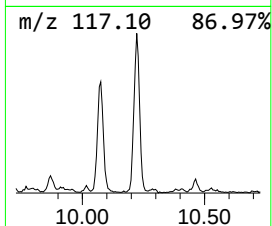
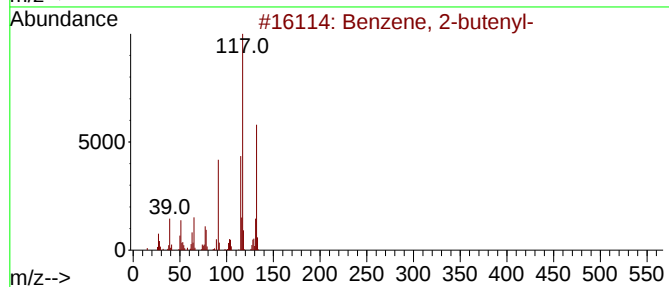
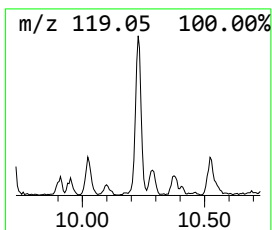
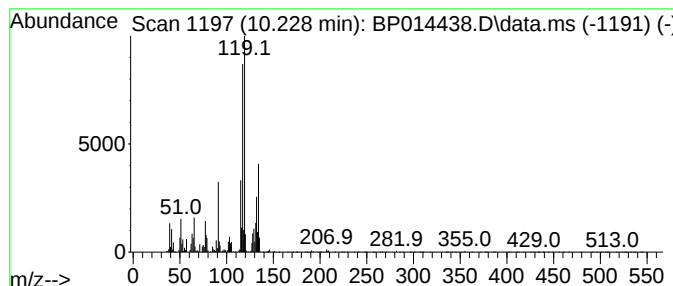
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 2-butenyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.228	2.58 ng/ul	217802	Naphthalene-d8	10.834

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	64
4			Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	64
5			Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	64



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
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Acq On : 31 Mar 2023 12:00
Operator : CG/JU
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ALS Vial : 47 Sample Multiplier: 1

Instrument :
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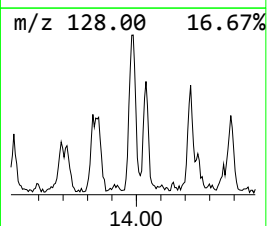
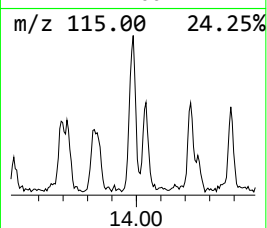
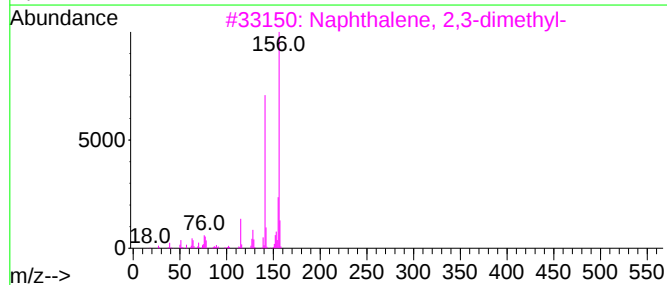
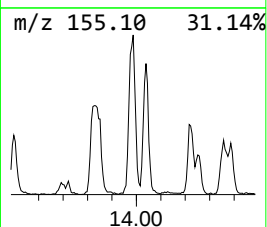
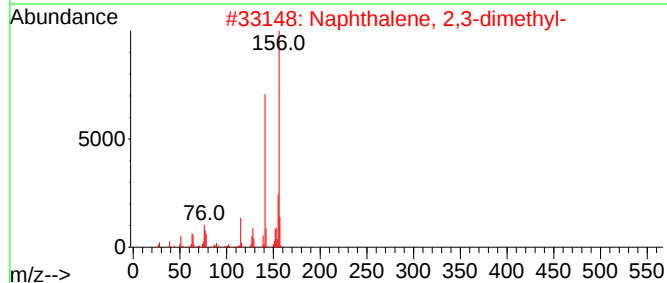
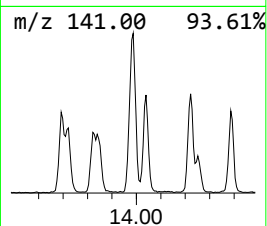
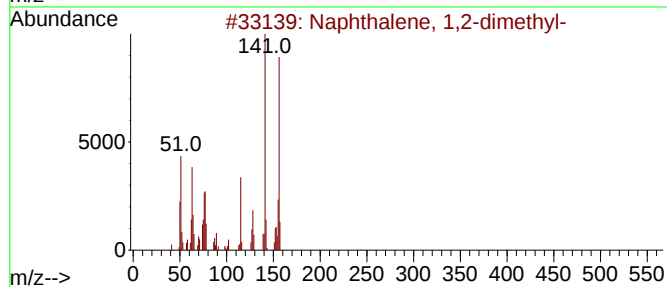
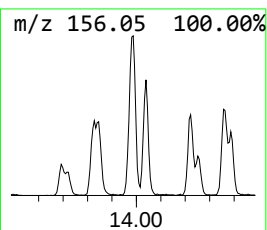
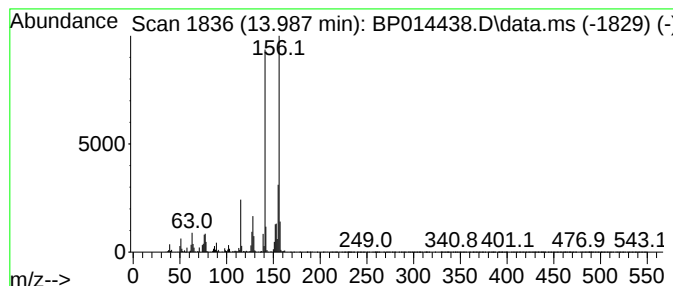
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Naphthalene, 1,2-dimethyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.987	2.84 ng/ul	362371	Acenaphthene-d10	14.663

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Naphthalene, 1,2-dimethyl-	156	C12H12	000573-98-8	97
2			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	97
3			Naphthalene, 2,3-dimethyl-	156	C12H12	000581-40-8	96
4			Naphthalene, 1,4-dimethyl-	156	C12H12	000571-58-4	96
5			Naphthalene, 1,7-dimethyl-	156	C12H12	000575-37-1	96



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP033023\
Data File : BP014438.D
Acq On : 31 Mar 2023 12:00
Operator : CG/JU
Sample : 02131-02DL 5X
Misc :
ALS Vial : 47 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP032823.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
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Benzene, 1,2,3-...	8.122	4.0	ng/ul	208583	1	8.010	1036360	20.0
Indane	8.387	2.3	ng/ul	117787	1	8.010	1036360	20.0
Indene	8.552	5.7	ng/ul	297261	1	8.010	1036360	20.0
Benzene, 1-ethy...	8.646	2.6	ng/ul	137454	1	8.010	1036360	20.0
Benzene, 2-ethy...	9.116	3.1	ng/ul	160372	1	8.010	1036360	20.0
Benzene, 2-bute...	10.228	2.6	ng/ul	217802	2	10.834	1691020	20.0
Naphthalene, 1,...	13.987	2.8	ng/ul	362371	3	14.663	2547910	20.0