

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
 Data File : BP022861.D
 Acq On : 05 Nov 2024 15:29
 Operator : RC/JU
 Sample : P4625-11
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AB6

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-BP100724.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP022861.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.181	29	33	36	rBV	13766	17967	0.58%	0.052%
2	3.605	101	105	112	rBV	23639	42822	1.37%	0.123%
3	3.899	150	155	164	rBV	33730	51355	1.65%	0.148%
4	4.399	234	240	247	rBV	8783	13050	0.42%	0.038%
5	5.340	388	400	409	rBV	47641	107948	3.46%	0.310%
6	5.681	453	458	464	rBV	9927	14803	0.48%	0.043%
7	6.146	531	537	544	rBV	29062	44563	1.43%	0.128%
8	6.622	613	618	624	rBV	7452	11748	0.38%	0.034%
9	6.787	638	646	648	rBV2	16417	24488	0.79%	0.070%
10	6.828	648	653	657	rBV	65761	115663	3.71%	0.333%
11	6.975	670	678	683	rBV	228460	367137	11.78%	1.056%
12	7.022	683	686	691	rVV	23216	33880	1.09%	0.097%
13	7.169	704	711	723	rBV	281045	481790	15.46%	1.386%
14	7.281	723	730	736	rVB	78967	125806	4.04%	0.362%
15	7.622	782	788	797	rBV	260044	414851	13.31%	1.193%
16	7.734	801	807	815	rVB2	23007	42516	1.36%	0.122%
17	7.987	842	850	861	rVV	1406750	2115883	67.91%	6.085%
18	8.146	873	877	885	rVV3	12970	24417	0.78%	0.070%
19	8.252	888	895	900	rVV3	7198	12697	0.41%	0.037%
20	8.334	900	909	927	rVV	259251	466049	14.96%	1.340%
21	8.705	962	972	977	rBV5	24972	57462	1.84%	0.165%
22	8.769	977	983	996	rVB2	333005	668611	21.46%	1.923%
23	8.957	1008	1015	1021	rVB	54671	94408	3.03%	0.271%
24	9.069	1023	1034	1041	rVV2	89630	211306	6.78%	0.608%
25	9.134	1041	1045	1055	rVV	60088	108049	3.47%	0.311%
26	9.222	1055	1060	1065	rVV	13461	22219	0.71%	0.064%
27	9.275	1065	1069	1075	rVB	9484	15212	0.49%	0.044%
28	9.475	1095	1103	1118	rBV	311805	522175	16.76%	1.502%
29	9.634	1124	1130	1141	rVB2	55939	100458	3.22%	0.289%
30	9.781	1144	1155	1166	rBV2	154544	283426	9.10%	0.815%
31	9.881	1167	1172	1176	rBV2	20954	36703	1.18%	0.106%
32	9.916	1176	1178	1184	rVB3	14041	19529	0.63%	0.056%
33	10.022	1188	1196	1209	rBV	456600	841479	27.01%	2.420%
34	10.381	1244	1257	1260	rVV	351532	624675	20.05%	1.796%
35	10.428	1260	1265	1272	rVV	1860704	3115945	100.00%	8.961%
36	10.546	1272	1285	1294	rVV2	326500	842240	27.03%	2.422%

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Integrator: RTE

Smoothing : OFF

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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-BP100724.MA.M
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37	10.622	1295	1298	1302	rVV2	20436	34838	1.12%	0.100%
38	10.669	1303	1306	1314	rVV2	11873	24376	0.78%	0.070%
39	10.793	1322	1327	1333	rVV2	23008	41314	1.33%	0.119%
40	11.034	1361	1368	1373	rVB4	8168	15483	0.50%	0.045%
41	11.146	1379	1387	1392	rBV3	7261	14437	0.46%	0.042%
42	11.240	1398	1403	1407	rVV8	5114	12670	0.41%	0.036%
43	11.281	1407	1410	1419	rVB2	7533	13492	0.43%	0.039%
44	11.481	1438	1444	1455	rBV	57197	113677	3.65%	0.327%
45	11.563	1455	1458	1467	rVB9	5462	12828	0.41%	0.037%
46	11.769	1486	1493	1494	rBV2	9057	16143	0.52%	0.046%
47	11.934	1515	1521	1529	rVV2	47064	83090	2.67%	0.239%
48	12.034	1532	1538	1545	rVV	115825	192963	6.19%	0.555%
49	12.163	1555	1560	1564	rVV2	10347	18971	0.61%	0.055%
50	12.257	1564	1576	1584	rVB2	229910	400287	12.85%	1.151%
51	12.381	1592	1597	1603	rVV5	16495	29340	0.94%	0.084%
52	12.651	1637	1643	1647	rVV4	11194	21699	0.70%	0.062%
53	12.698	1647	1651	1655	rVV6	7877	15406	0.49%	0.044%
54	12.745	1657	1659	1667	rVV7	7525	13236	0.42%	0.038%
55	12.987	1694	1700	1705	rBV7	7184	16066	0.52%	0.046%
56	13.257	1742	1746	1748	rVV4	11290	17481	0.56%	0.050%
57	13.281	1748	1750	1757	rVB7	15318	25294	0.81%	0.073%
58	13.416	1761	1773	1782	rBV7	13780	51360	1.65%	0.148%
59	13.563	1789	1798	1803	rBV3	20672	46049	1.48%	0.132%
60	13.651	1806	1813	1825	rVV	1021783	1356632	43.54%	3.901%
61	13.792	1833	1837	1840	rBV2	18282	26731	0.86%	0.077%
62	13.828	1840	1843	1850	rVB5	13136	24457	0.78%	0.070%
63	13.934	1854	1861	1877	rVB	1097430	1532179	49.17%	4.406%
64	14.057	1877	1882	1888	rVB9	6302	12740	0.41%	0.037%
65	14.245	1899	1914	1919	rBV	645803	939675	30.16%	2.702%
66	14.304	1919	1924	1932	rVV	498638	758839	24.35%	2.182%
67	14.387	1933	1938	1945	rVB	13648	24740	0.79%	0.071%
68	14.516	1953	1960	1972	rBV3	49497	161912	5.20%	0.466%
69	14.657	1980	1984	1989	rVB	12673	19656	0.63%	0.057%
70	14.745	1995	1999	2005	rVB	59477	83428	2.68%	0.240%
71	15.116	2058	2062	2072	rVB	33190	58426	1.88%	0.168%
72	15.257	2079	2086	2092	rBV	1248487	1993311	63.97%	5.732%
73	15.322	2092	2097	2105	rVB	169507	256692	8.24%	0.738%
74	15.398	2105	2110	2119	rBV	408340	563257	18.08%	1.620%
75	15.475	2119	2123	2129	rVB7	16017	29693	0.95%	0.085%
76	15.639	2145	2151	2160	rVB	64396	108351	3.48%	0.312%

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Integrator: RTE
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 Start Thrs: 0.2 Max Peaks: 100
 Stop Thrs : 0 Peak Location: TOP

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77	15.722	2160	2165	2169	rBV3	14503	29943	0.96%	0.086%
78	15.757	2169	2171	2176	rVB3	11749	13639	0.44%	0.039%
79	15.992	2205	2211	2215	rBV8	9870	24374	0.78%	0.070%
80	16.204	2242	2247	2252	rVV3	18159	28661	0.92%	0.082%
81	16.316	2261	2266	2272	rBV4	12252	24134	0.77%	0.069%
82	16.851	2345	2357	2360	rBV6	12530	37522	1.20%	0.108%
83	16.957	2372	2375	2382	rVB6	8161	14158	0.45%	0.041%
84	17.051	2383	2391	2398	rBV	774148	1228422	39.42%	3.533%
85	17.157	2402	2409	2418	rVV2	1550914	2347177	75.33%	6.750%
86	17.486	2459	2465	2472	rBV	164122	241970	7.77%	0.696%
87	17.851	2522	2527	2532	rBV5	11644	19943	0.64%	0.057%
88	17.933	2537	2541	2547	rVB	26923	39884	1.28%	0.115%
89	17.998	2547	2552	2557	rBV3	52020	105036	3.37%	0.302%
90	18.104	2567	2570	2577	rVB	15713	26901	0.86%	0.077%
91	18.286	2595	2601	2604	rBV6	18297	35338	1.13%	0.102%
92	18.510	2634	2639	2640	rBV5	8239	12174	0.39%	0.035%
93	19.075	2731	2735	2740	rVB3	25586	33573	1.08%	0.097%
94	19.157	2745	2749	2753	rBV	18782	29038	0.93%	0.084%
95	19.327	2774	2778	2784	rBV8	15055	28256	0.91%	0.081%
96	19.533	2806	2813	2821	rBV	2071554	2965068	95.16%	8.527%
97	20.157	2916	2919	2924	rBV2	37746	60825	1.95%	0.175%
98	21.498	3141	3147	3157	rVB2	945103	1497431	48.06%	4.306%
99	24.527	3653	3662	3675	rVB	1137400	3109752	99.80%	8.943%
100	24.745	3692	3699	3712	rVB	543226	1545647	49.60%	4.445%

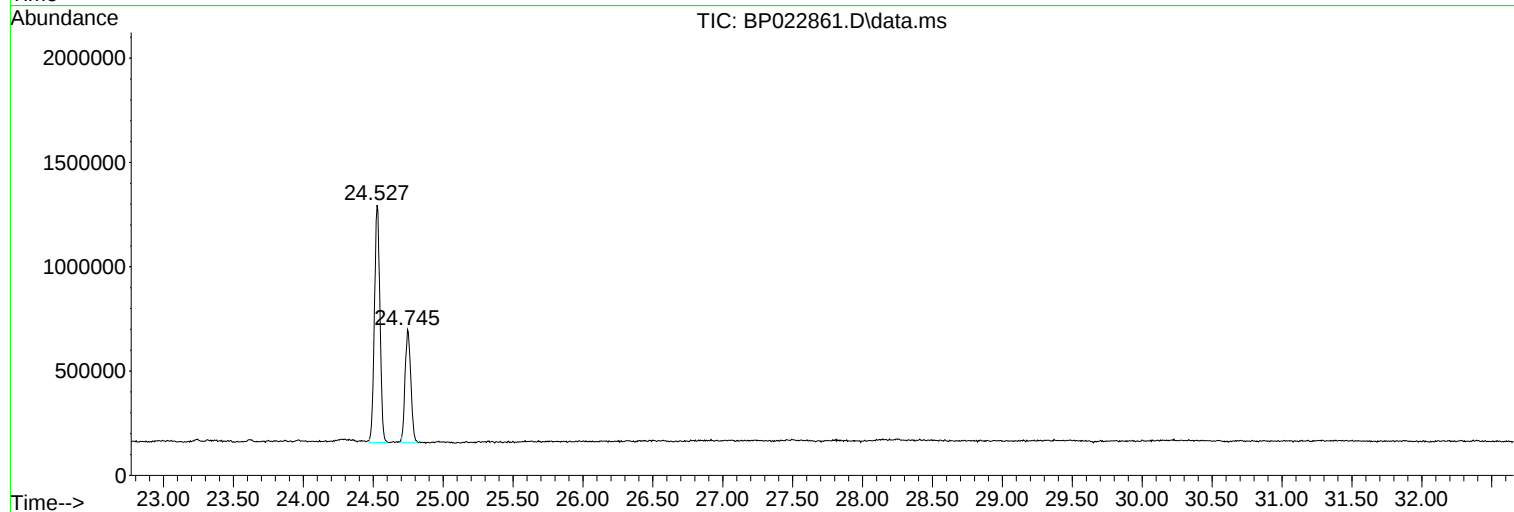
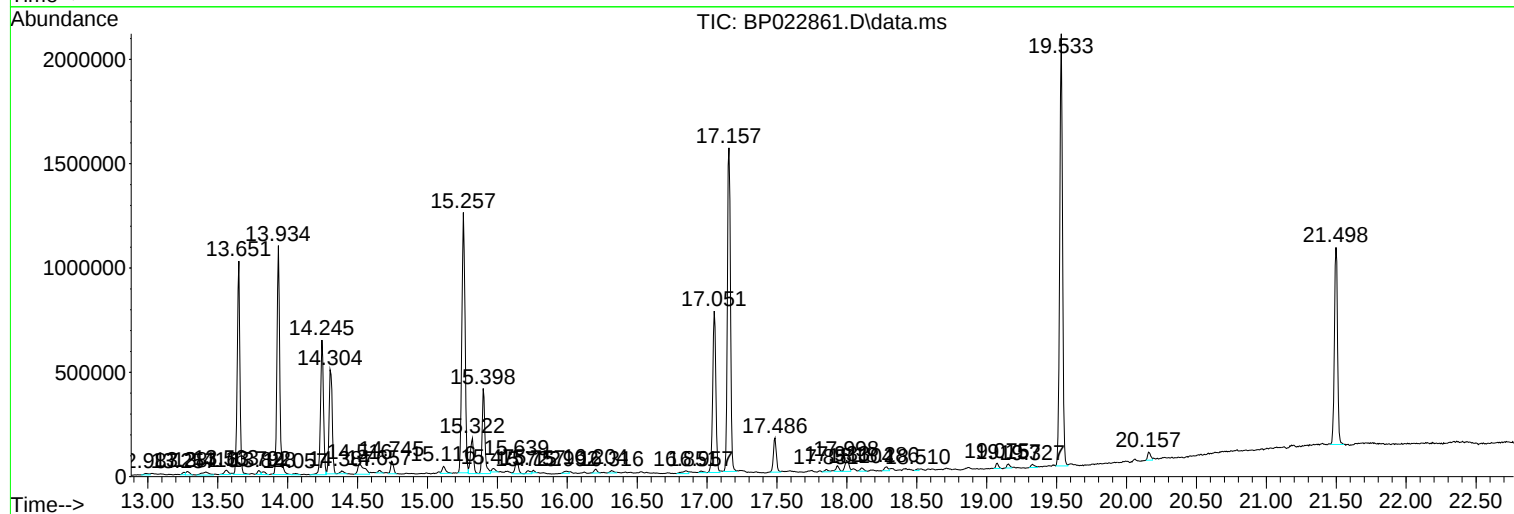
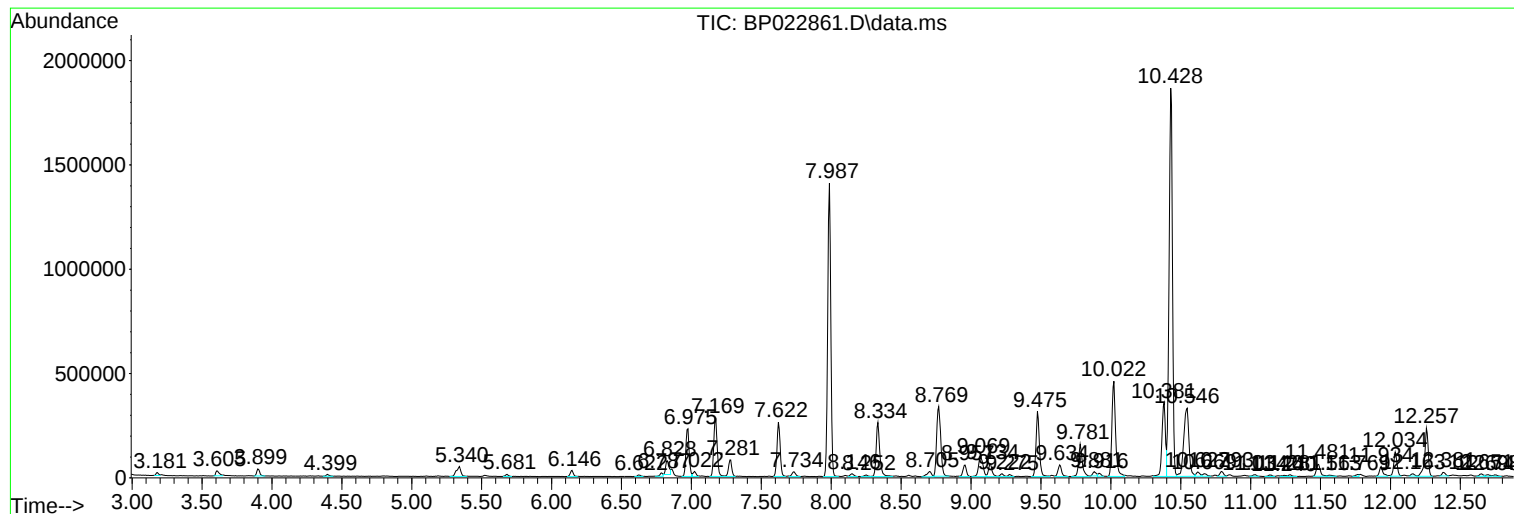
Sum of corrected areas: 34773415

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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



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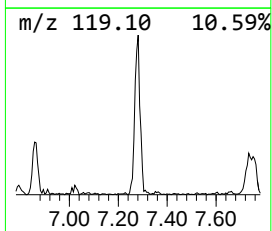
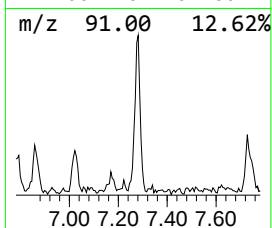
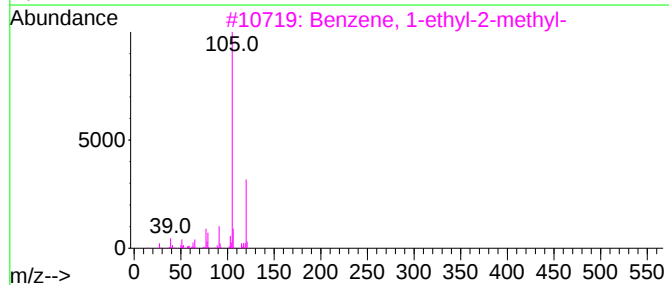
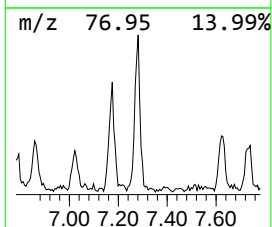
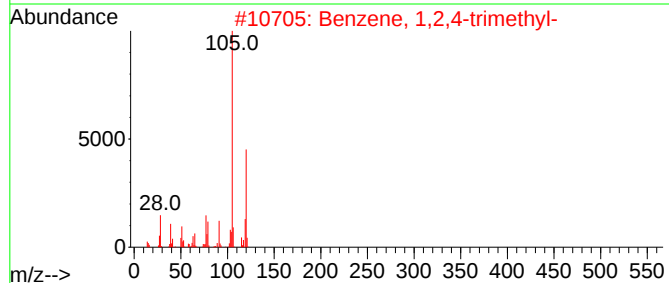
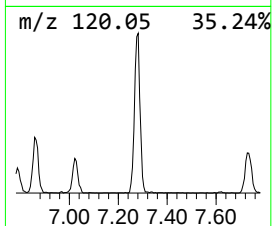
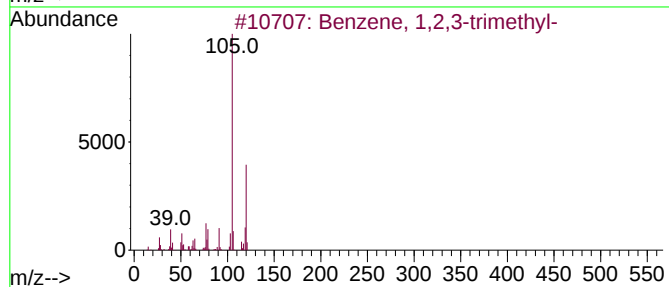
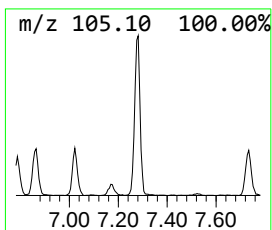
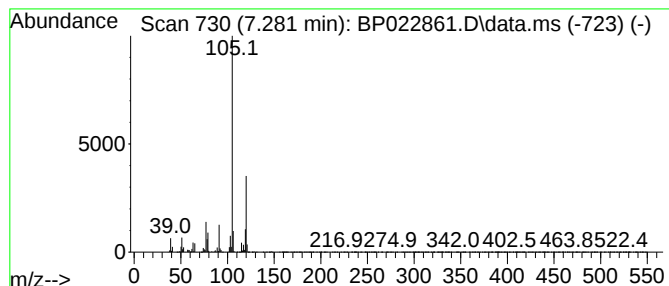
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 4 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.281	6.07 ng/ul	125806	1,4-Dichlorobenzene-d4	7.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	97
2			Benzene, 1,2,4-trimethyl-	120	C9H12	000095-63-6	94
3			Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	93
4			Mesitylene	120	C9H12	000108-67-8	91
5			Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91



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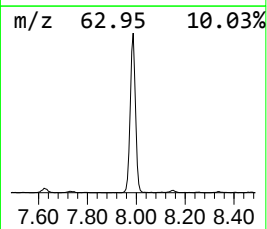
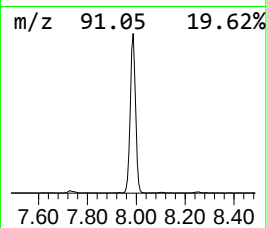
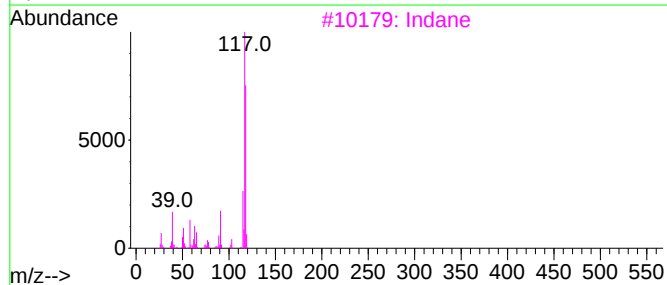
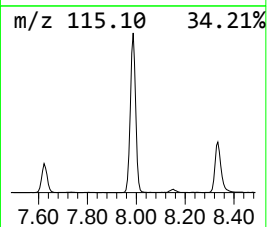
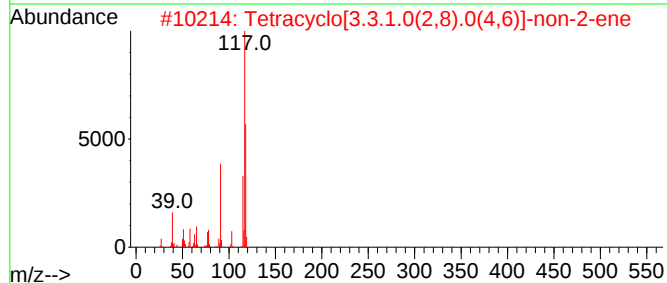
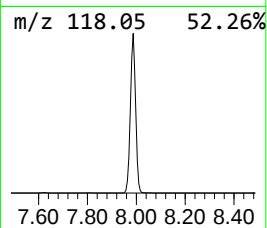
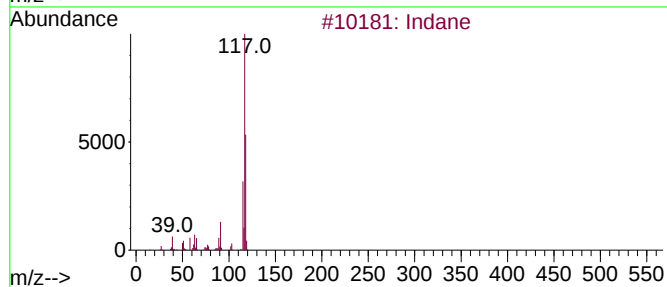
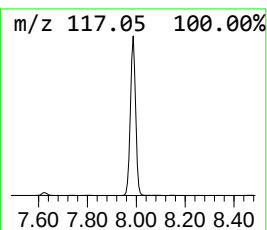
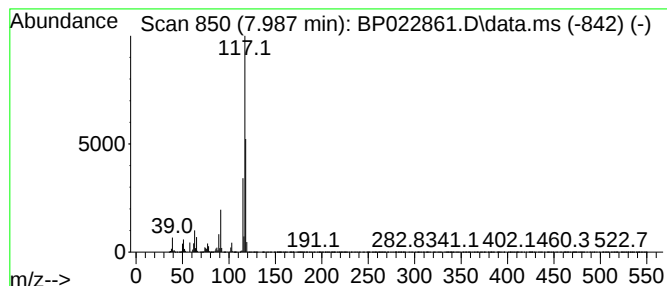
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.987	102.01 ng/ul	2115880	1,4-Dichlorobenzene-d4	7.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Indane	118	C9H10	000496-11-7	94
2			Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118	C9H10	1000191-13-7	72
3			Indane	118	C9H10	000496-11-7	64
4			Benzene, cyclopropyl-	118	C9H10	000873-49-4	64
5			1,3-Methanopentalene, 1,2,3,5-te...	118	C9H10	128600-88-4	59



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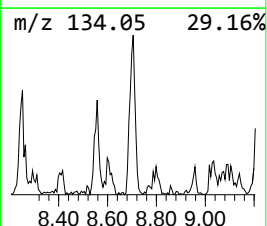
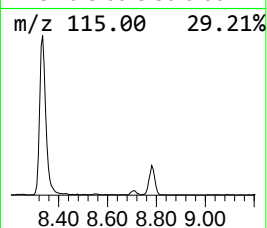
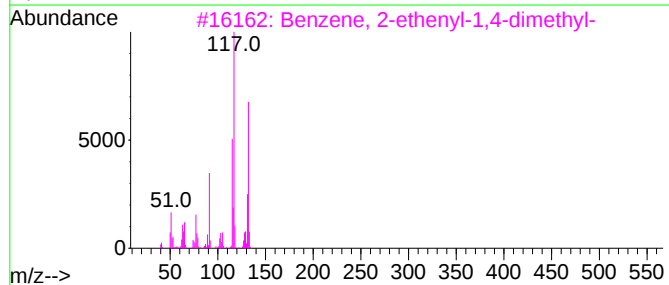
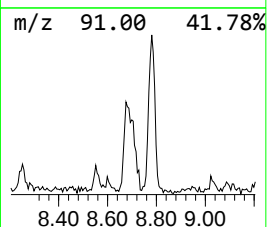
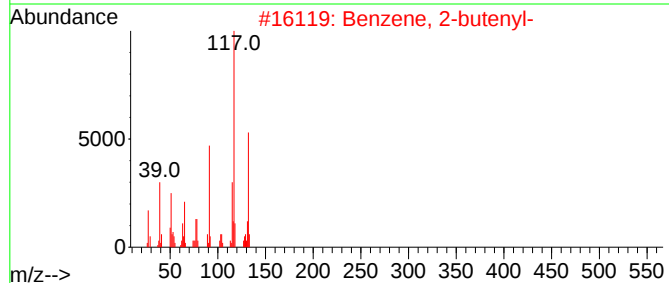
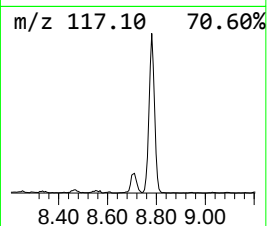
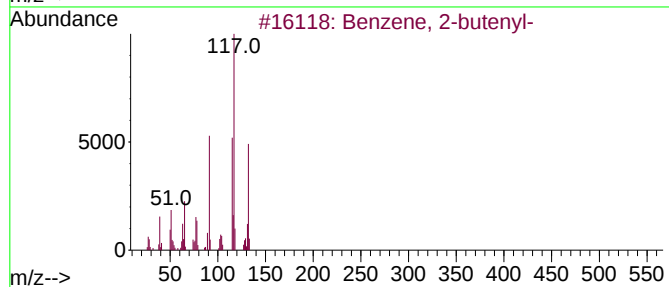
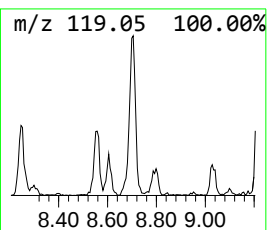
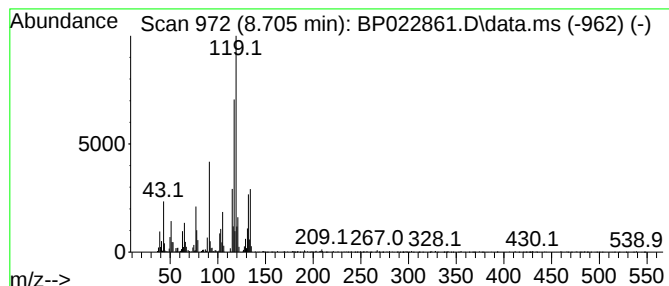
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 2-butenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.705	2.77 ng/ul	57462	1,4-Dichlorobenzene-d4	7.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-butenyl-	132	C10H12	001560-06-1	93
2			Benzene, 2-butenyl-	132	C10H12	001560-06-1	80
3			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	56
4			p-Cymene	134	C10H14	000099-87-6	55
5			o-Cymene	134	C10H14	000527-84-4	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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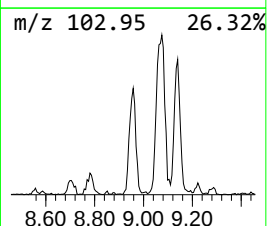
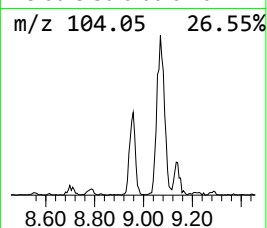
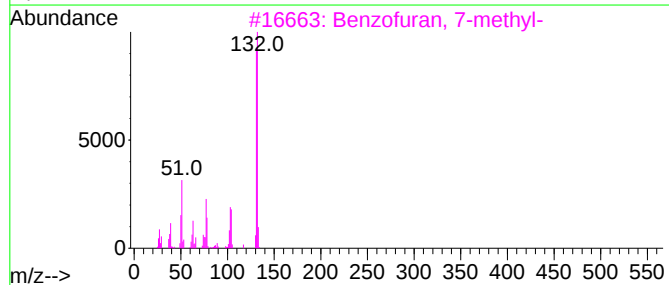
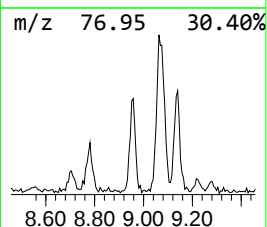
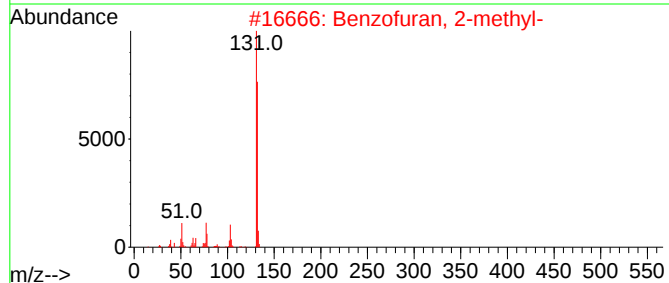
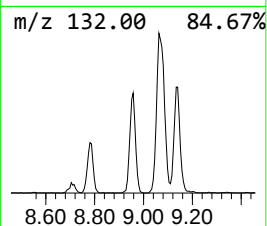
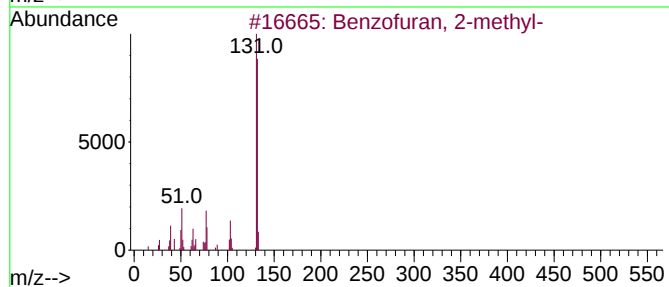
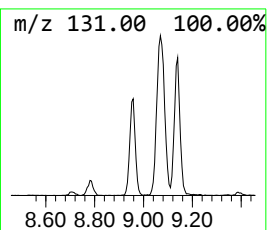
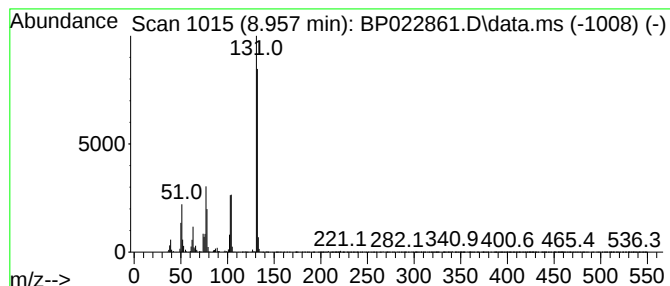
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 Benzofuran, 2-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.957	4.55 ng/ul	94408	1,4-Dichlorobenzene-d4	7.622

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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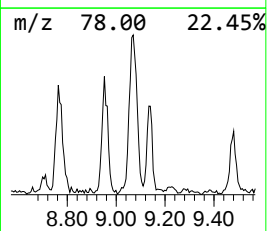
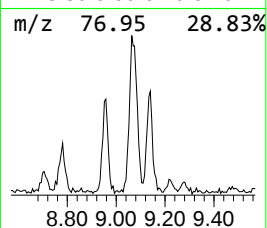
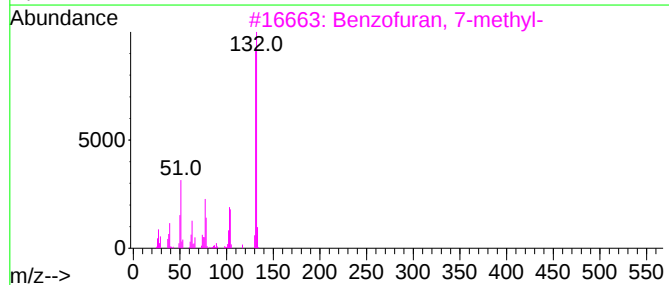
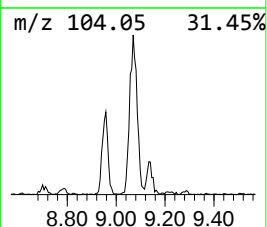
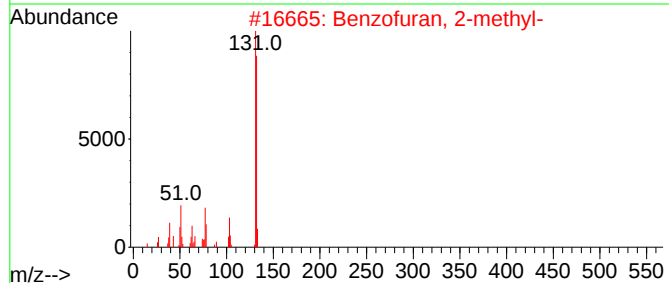
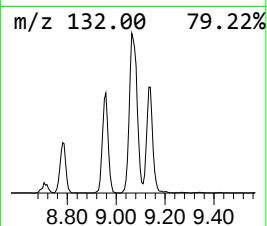
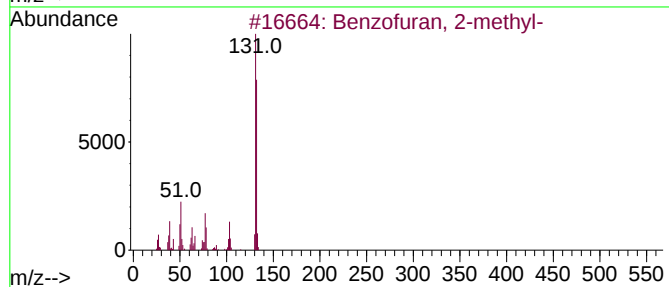
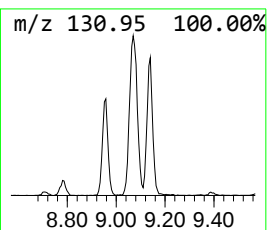
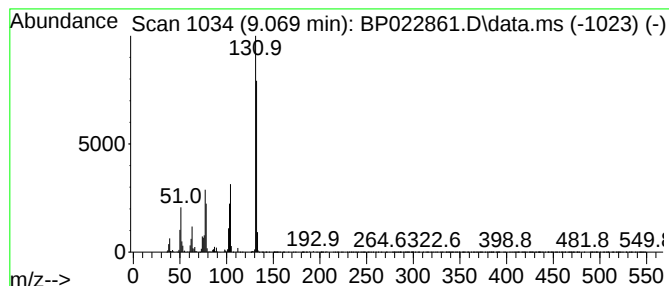
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzfuran, 7-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.069	6.77 ng/ul	211306	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3			Benzofuran, 7-methyl-	132	C9H8O	017059-52-8	94
4			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	93
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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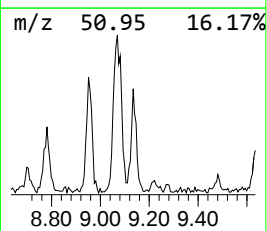
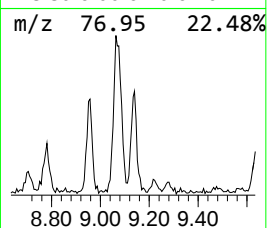
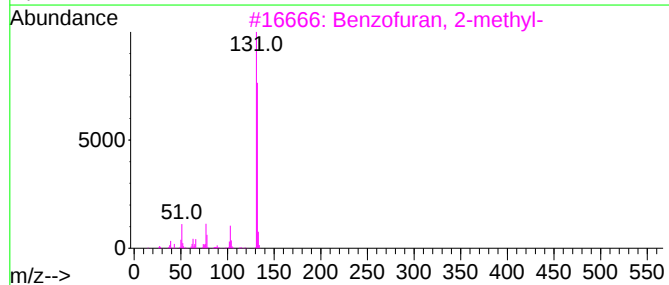
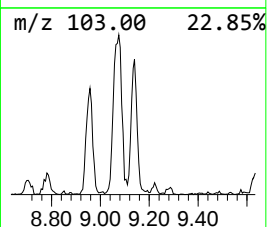
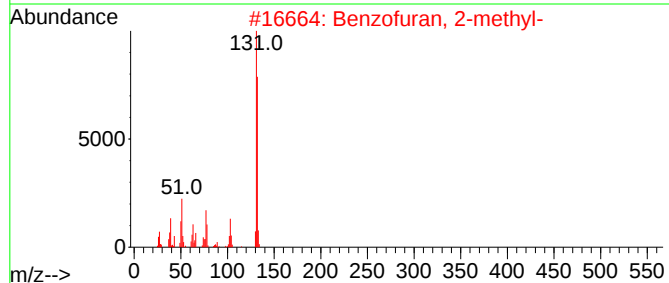
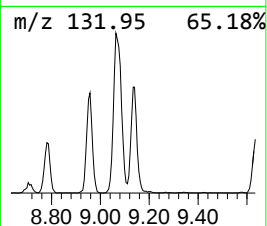
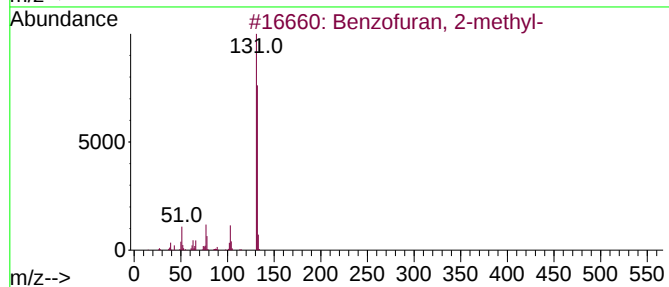
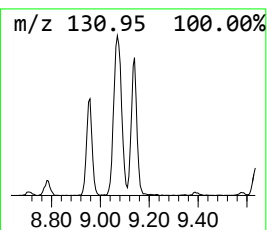
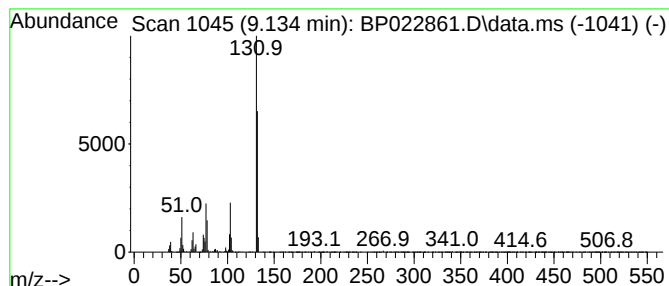
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 Cinnamaldehyde, (E)- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.134	3.46 ng/ul	108049	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
2			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	94
3			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	91
4			Cinnamaldehyde, (E)-	132	C9H8O	014371-10-9	91
5			Benzofuran, 2-methyl-	132	C9H8O	004265-25-2	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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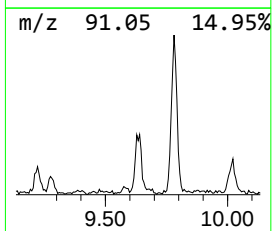
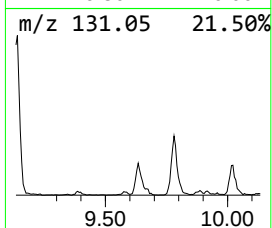
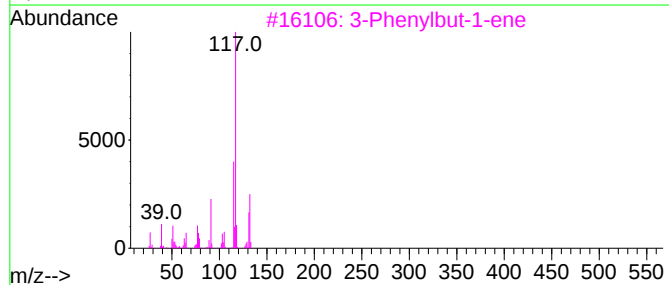
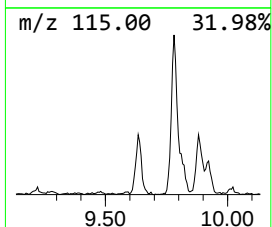
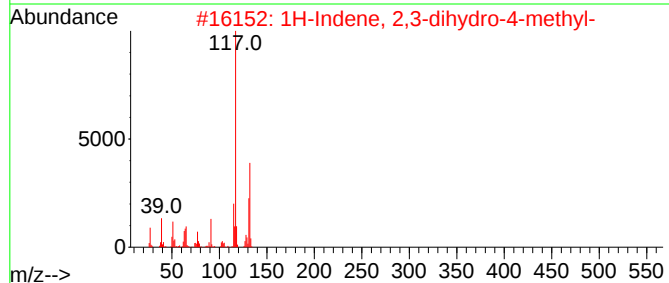
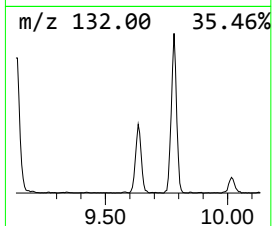
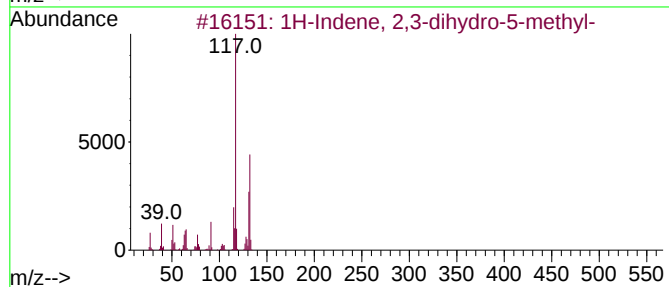
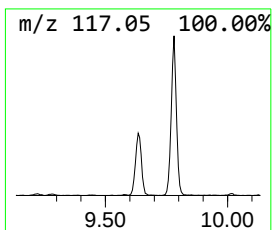
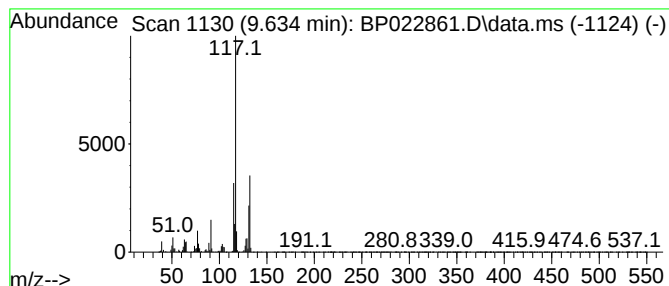
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 1H-Indene, 2,3-dihydro-5-me... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.634	3.22 ng/ul	100458	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	91
2			1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	90
3			3-Phenylbut-1-ene	132	C10H12	000934-10-1	87
4			Indan, 1-methyl-	132	C10H12	000767-58-8	83
5			Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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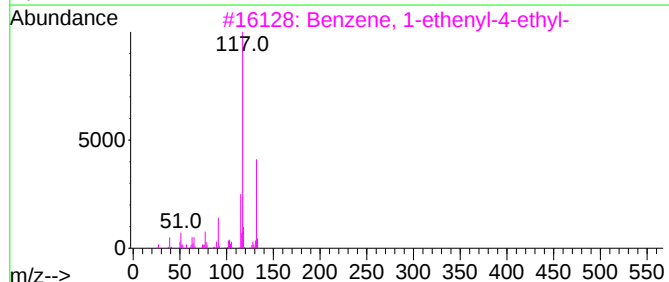
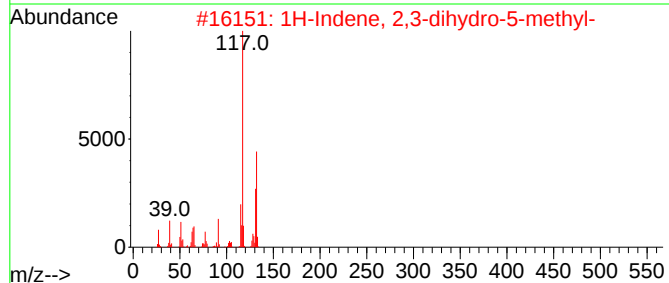
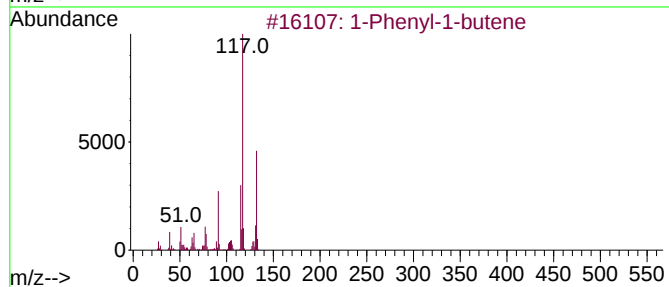
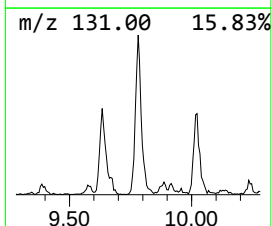
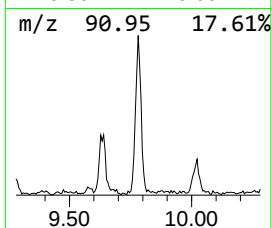
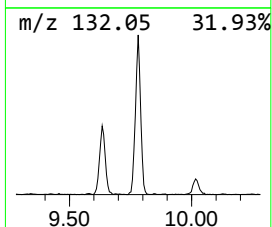
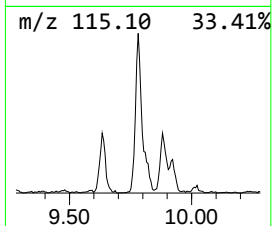
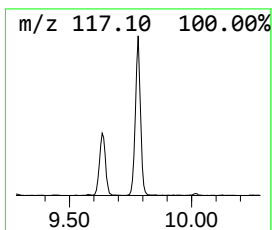
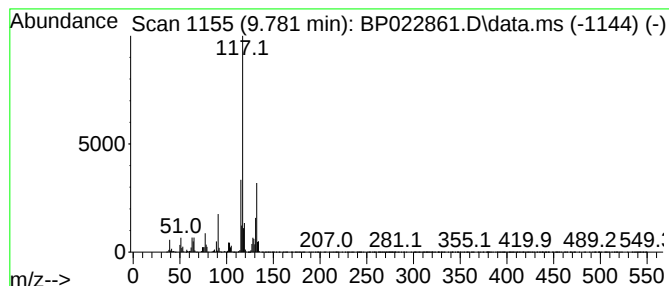
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Quant Title : SVOA CALIBRATION

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

Peak Number 12 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.781	9.07 ng/ul	283426	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1-Phenyl-1-butene	132	C10H12	000824-90-8	90
2			1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	87
3			Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	87
4			Indan, 1-methyl-	132	C10H12	000767-58-8	87
5			Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
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Instrument :
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ClientSampleId :
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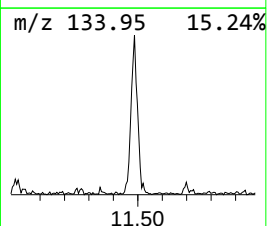
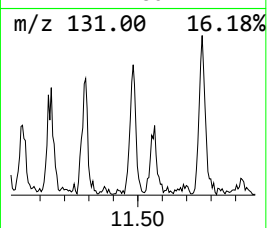
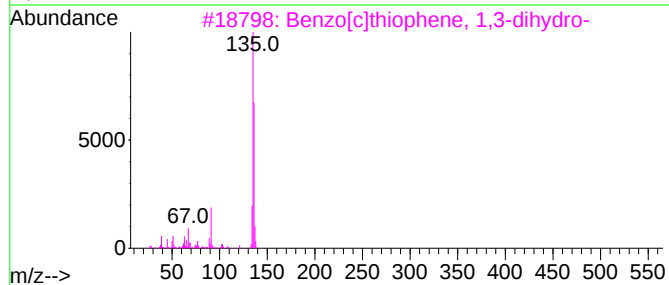
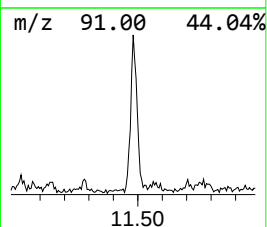
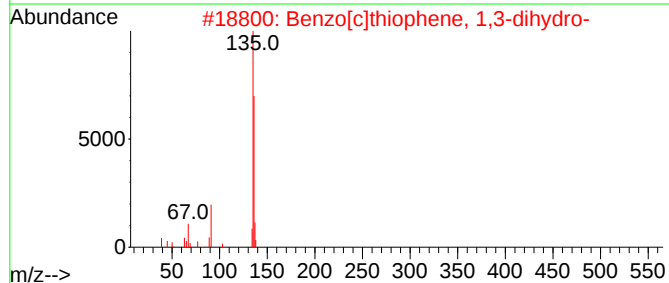
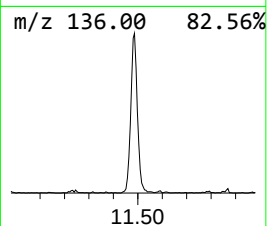
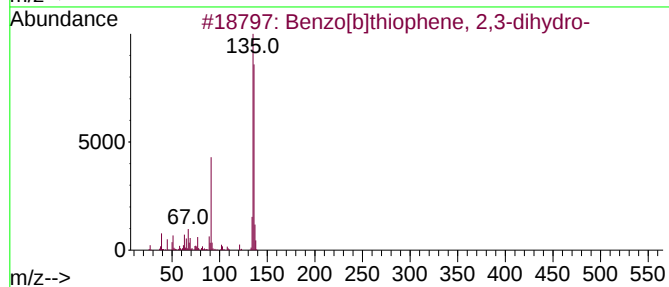
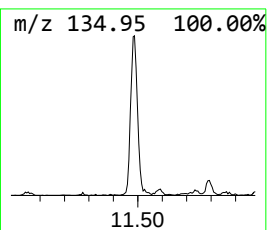
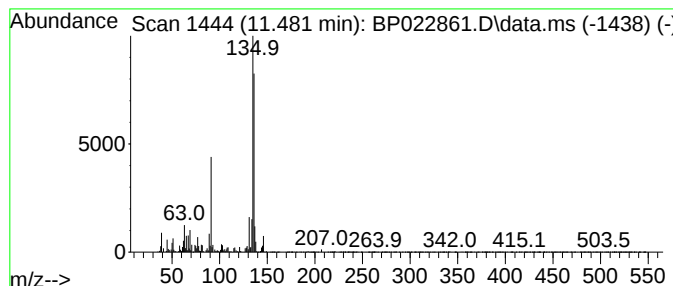
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Quant Title : SVOA CALIBRATION

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

Peak Number 13 Benzo[b]thiophene, 2,3-dihy... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.481	3.64 ng/ul	113677	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	93
2			Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	90
3			Benzo[c]thiophene, 1,3-dihydro-	136	C8H8S	002471-92-3	87
4			4-Hydroxy-3-methylbenzaldehyde	136	C8H8O2	015174-69-3	72
5			Benzo[b]thiophene, 2,3-dihydro-	136	C8H8S	004565-32-6	70



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
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Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

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ClientSampleId :
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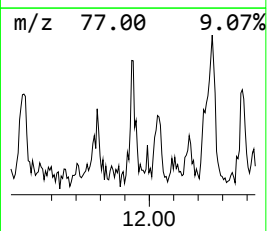
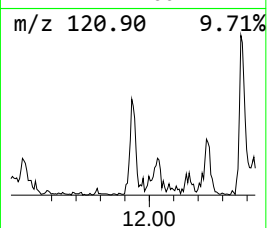
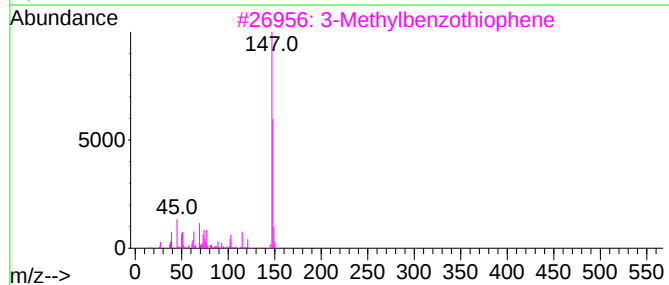
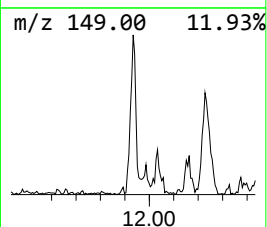
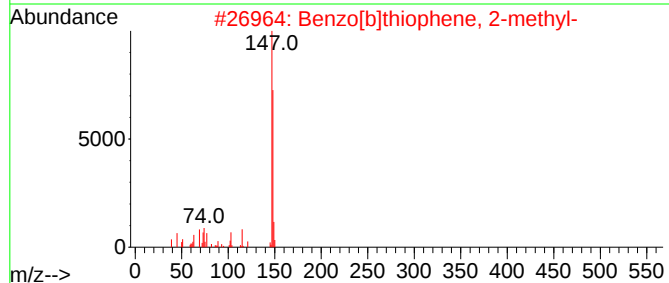
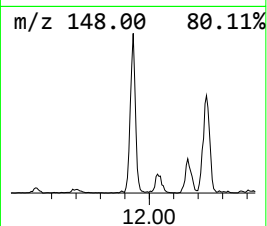
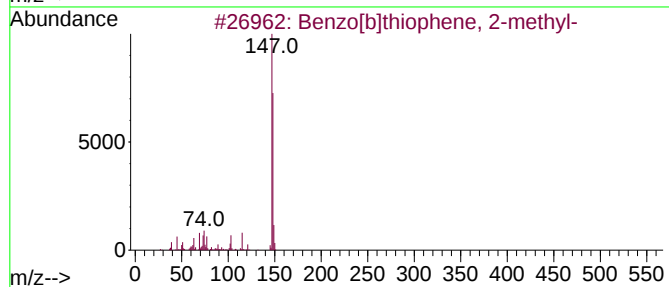
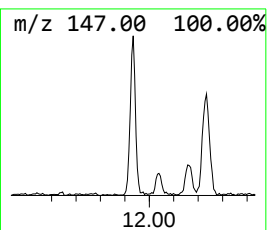
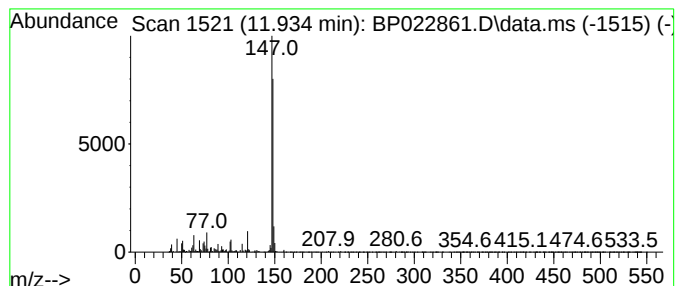
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Quant Title : SVOA CALIBRATION

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

Peak Number 14 Benzo[b]thiophene, 2-methyl- Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.934	2.66 ng/ul	83090	Naphthalene-d8	10.381

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
2			Benzo[b]thiophene, 2-methyl-	148	C9H8S	001195-14-8	91
3			3-Methylbenzothiophene	148	C9H8S	001455-18-1	90
4			Benzo[b]thiophene, 4-methyl-	148	C9H8S	014315-11-8	87
5			Benzo[b]thiophene, 6-methyl-	148	C9H8S	016587-47-6	86



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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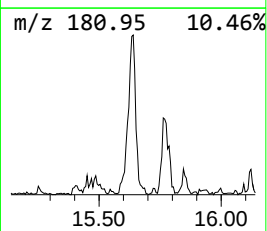
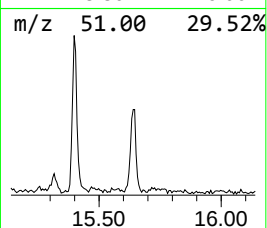
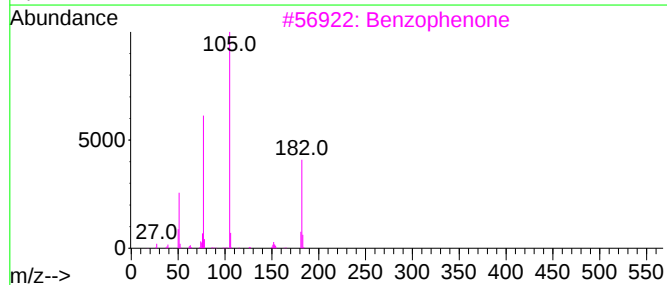
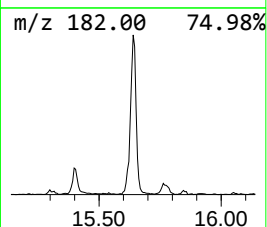
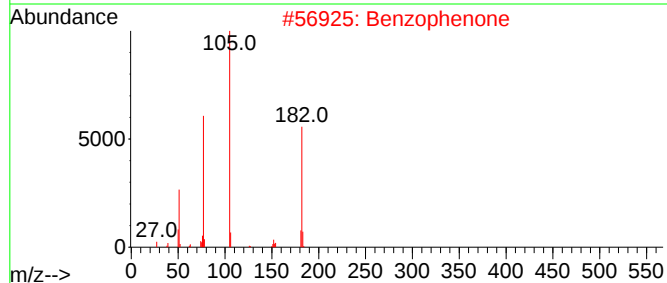
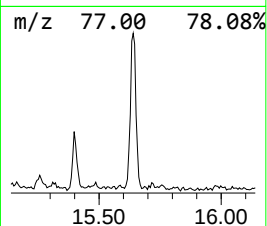
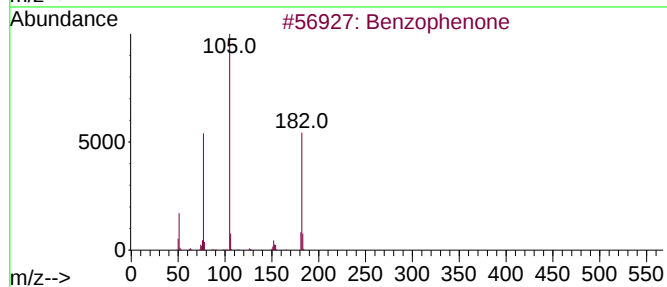
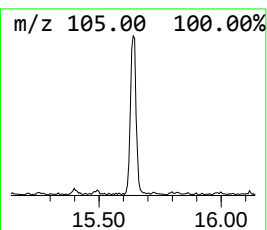
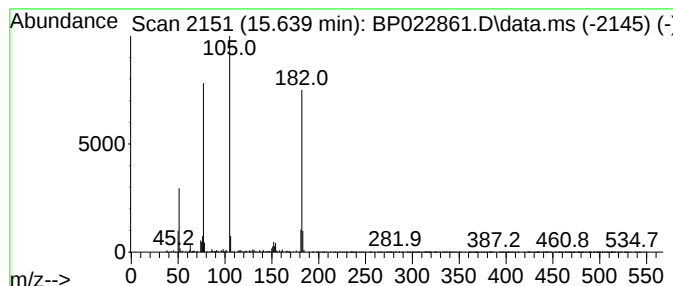
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Quant Title : SVOA CALIBRATION

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

Peak Number 15 Benzophenone Concentration Rank 13

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.640	2.31 ng/ul	108351	Acenaphthene-d10	14.245

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzophenone	182	C13H10O	000119-61-9	97
2			Benzophenone	182	C13H10O	000119-61-9	96
3			Benzophenone	182	C13H10O	000119-61-9	94
4			Benzophenone	182	C13H10O	000119-61-9	91
5			Benzophenone	182	C13H10O	000119-61-9	90



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP110424\
Data File : BP022861.D
Acq On : 05 Nov 2024 15:29
Operator : RC/JU
Sample : P4625-11
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
C0AB6

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP100724.MA.M
Quant Title : SVOA CALIBRATION

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

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1,2,3-...	7.281	6.1	ng/ul	125806	1	7.622	414851	20.0
Indane	7.987	102.0	ng/ul	2115880	1	7.622	414851	20.0
Benzene, 2-bute...	8.705	2.8	ng/ul	57462	1	7.622	414851	20.0
Benzofuran, 2-m...	8.957	4.5	ng/ul	94408	1	7.622	414851	20.0
Benzofuran, 7-m...	9.069	6.8	ng/ul	211306	2	10.381	624675	20.0
Cinnamaldehyde,...	9.134	3.5	ng/ul	108049	2	10.381	624675	20.0
1H-Indene, 2,3-...	9.634	3.2	ng/ul	100458	2	10.381	624675	20.0
1-Phenyl-1-butene	9.781	9.1	ng/ul	283426	2	10.381	624675	20.0
Benzo[b]thiophe...	11.481	3.6	ng/ul	113677	2	10.381	624675	20.0
Benzo[b]thiophe...	11.934	2.7	ng/ul	83090	2	10.381	624675	20.0
Benzophenone	15.640	2.3	ng/ul	108351	3	14.245	939675	20.0