

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
 Data File : BP010892.D
 Acq On : 27 Jun 2022 23:53
 Operator : CG/JU
 Sample : N3401-14DL 5X
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 C0AF2DL

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-BP062322.M
 Title : SVOA CALIBRATION

Signal : TIC: BP010892.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.452	58	62	73	rVB2	122863	193535	3.42%	0.256%
2	3.617	87	90	100	rBV	194318	276762	4.89%	0.367%
3	3.928	138	143	148	rVB	46622	64465	1.14%	0.085%
4	4.099	167	172	176	rBV2	28307	41487	0.73%	0.055%
5	4.234	186	195	203	rBV5	17157	36547	0.65%	0.048%
6	4.375	214	219	226	rVB4	37646	81378	1.44%	0.108%
7	4.464	230	234	240	rVB5	20912	34797	0.62%	0.046%
8	4.864	298	302	308	rVV3	43805	79991	1.41%	0.106%
9	4.922	309	312	320	rVV6	16479	35951	0.64%	0.048%
10	5.228	359	364	372	rVB	141273	206392	3.65%	0.273%
11	5.381	385	390	394	rBV	64096	94631	1.67%	0.125%
12	5.422	394	397	406	rVB4	31287	55388	0.98%	0.073%
13	6.205	521	530	538	rBV	1333079	1973267	34.88%	2.614%
14	6.693	606	613	621	rVB	2093382	3104822	54.88%	4.113%
15	6.881	635	645	656	rVB	129008	234207	4.14%	0.310%
16	7.052	666	674	679	rBV	462332	709942	12.55%	0.941%
17	7.099	679	682	685	rVV2	25395	39747	0.70%	0.053%
18	7.134	685	688	694	rVB2	19413	35758	0.63%	0.047%
19	7.240	699	706	713	rBV	476038	722399	12.77%	0.957%
20	7.357	720	726	732	rVB	401870	591781	10.46%	0.784%
21	7.575	758	763	766	rVV	50391	80983	1.43%	0.107%
22	7.610	766	769	774	rVV	64475	90433	1.60%	0.120%
23	7.705	776	785	793	rVV	1999462	3184324	56.29%	4.219%
24	7.822	799	805	814	rVB	222328	357834	6.33%	0.474%
25	8.016	830	838	842	rVV	27696	49350	0.87%	0.065%
26	8.075	842	848	858	rVV	1347130	2121817	37.51%	2.811%
27	8.199	862	869	875	rVV	314035	451750	7.99%	0.598%
28	8.263	875	880	887	rVB	41038	61167	1.08%	0.081%
29	8.352	887	895	899	rBV2	291824	456907	8.08%	0.605%
30	8.416	899	906	914	rVV2	340535	670420	11.85%	0.888%
31	8.510	917	922	932	rVB	47373	92526	1.64%	0.123%
32	8.657	942	947	952	rBV	88378	132816	2.35%	0.176%
33	8.710	952	956	963	rVV	49841	77720	1.37%	0.103%
34	8.810	966	973	978	rVV	1890609	2825380	49.94%	3.743%
35	8.869	978	983	995	rVB3	519123	1435761	25.38%	1.902%
36	9.046	1008	1013	1020	rBV2	41774	74726	1.32%	0.099%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
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37	9.140	1023	1029	1035	rBV2	44189	79063	1.40%	0.105%
38	9.293	1047	1055	1056	rBV2	33018	47480	0.84%	0.063%
39	9.334	1056	1062	1067	rVV	1264135	1906126	33.69%	2.525%
40	9.393	1067	1072	1078	rVB	293536	454583	8.04%	0.602%

41	9.587	1098	1105	1117	rBV2	332023	594432	10.51%	0.788%
42	9.704	1119	1125	1127	rBV3	33959	53652	0.95%	0.071%
43	9.752	1127	1133	1145	rVB	623697	1009692	17.85%	1.338%
44	9.904	1148	1159	1167	rBV3	1989023	3493372	61.75%	4.628%
45	9.981	1167	1172	1176	rBV2	69118	123684	2.19%	0.164%

46	10.122	1186	1196	1205	rVB2	568838	1062469	18.78%	1.408%
47	10.210	1205	1211	1219	rVB	107811	181136	3.20%	0.240%
48	10.375	1235	1239	1245	rBV4	13041	28621	0.51%	0.038%
49	10.428	1245	1248	1252	rBV2	19691	33126	0.59%	0.044%
50	10.504	1252	1261	1265	rVV	2726019	4676415	82.67%	6.195%

51	10.551	1265	1269	1277	rVV	3533404	5657061	100.00%	7.495%
52	10.646	1277	1285	1293	rVV	435927	749856	13.26%	0.993%
53	10.722	1293	1298	1304	rVV	91978	135451	2.39%	0.179%
54	10.787	1304	1309	1315	rVB	42801	69656	1.23%	0.092%
55	10.875	1318	1324	1329	rBV3	31363	60285	1.07%	0.080%

56	11.075	1351	1358	1362	rBV9	15096	33223	0.59%	0.044%
57	11.134	1362	1368	1370	rVV	33326	58697	1.04%	0.078%
58	11.163	1370	1373	1382	rVB4	35517	71800	1.27%	0.095%
59	11.340	1395	1403	1407	rBV	24079	40663	0.72%	0.054%
60	11.422	1412	1417	1422	rVB	40793	64171	1.13%	0.085%

61	11.616	1446	1450	1457	rVV3	29990	52813	0.93%	0.070%
62	11.698	1459	1464	1470	rVV4	17729	31008	0.55%	0.041%
63	11.834	1480	1487	1492	rVV	49280	87341	1.54%	0.116%
64	11.898	1492	1498	1506	rVB4	39569	79644	1.41%	0.106%
65	12.169	1535	1544	1551	rVB2	66941	131876	2.33%	0.175%

66	12.240	1551	1556	1560	rBV	31625	49141	0.87%	0.065%
67	12.398	1575	1583	1590	rVB5	78120	134964	2.39%	0.179%
68	12.481	1592	1597	1604	rVB3	43534	74109	1.31%	0.098%
69	12.710	1628	1636	1642	rBV2	21179	48538	0.86%	0.064%
70	12.816	1649	1654	1660	rVV	169982	255116	4.51%	0.338%

71	12.887	1663	1666	1672	rVB8	17763	33317	0.59%	0.044%
72	13.069	1693	1697	1706	rBV4	19058	34743	0.61%	0.046%
73	13.328	1736	1741	1744	rBV2	92261	135822	2.40%	0.180%
74	13.363	1744	1747	1758	rVB4	46013	79622	1.41%	0.105%
75	13.492	1764	1769	1773	rBV3	94366	154467	2.73%	0.205%

76	13.534	1773	1776	1779	rVV	44050	66368	1.17%	0.088%
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 Title : SVOA CALIBRATION

77	13.575	1779	1783	1789	rVV	88801	130600	2.31%	0.173%
78	13.663	1794	1798	1806	rVB7	18567	34247	0.61%	0.045%
79	13.781	1812	1818	1825	rVB	1035863	1379869	24.39%	1.828%
80	13.981	1845	1852	1857	rBV3	106086	176455	3.12%	0.234%
81	14.051	1858	1864	1877	rVV	1081918	1613174	28.52%	2.137%
82	14.187	1877	1887	1895	rVB4	41530	97360	1.72%	0.129%
83	14.281	1898	1903	1909	rVB6	26604	40766	0.72%	0.054%
84	14.363	1909	1917	1930	rBV	3814662	5350597	94.58%	7.089%
85	14.522	1937	1944	1946	rBV	40060	67167	1.19%	0.089%
86	14.575	1949	1953	1959	rVB	53373	91329	1.61%	0.121%
87	14.692	1968	1973	1976	rBV3	21434	31586	0.56%	0.042%
88	14.851	1994	2000	2002	rBV4	16265	31787	0.56%	0.042%
89	14.887	2002	2006	2014	rVB	38972	59286	1.05%	0.079%
90	15.069	2032	2037	2049	rVB	41725	73408	1.30%	0.097%
91	15.263	2063	2070	2075	rVB2	23838	45676	0.81%	0.061%
92	15.363	2079	2087	2100	rBV	1525731	2178954	38.52%	2.887%
93	15.481	2102	2107	2113	rBV	192506	288338	5.10%	0.382%
94	15.604	2123	2128	2136	rVB2	43820	60922	1.08%	0.081%
95	17.128	2380	2387	2397	rBV	3858044	5395434	95.38%	7.148%
96	17.228	2397	2404	2419	rVB	1375539	2011530	35.56%	2.665%
97	19.475	2781	2786	2792	rBV	1682555	2142090	37.87%	2.838%
98	21.239	3081	3086	3092	rBV	3972595	4715456	83.36%	6.247%
99	23.445	3456	3461	3470	rVB	944687	1811688	32.03%	2.400%
100	23.604	3482	3488	3499	rVB2	2480099	4742351	83.83%	6.283%

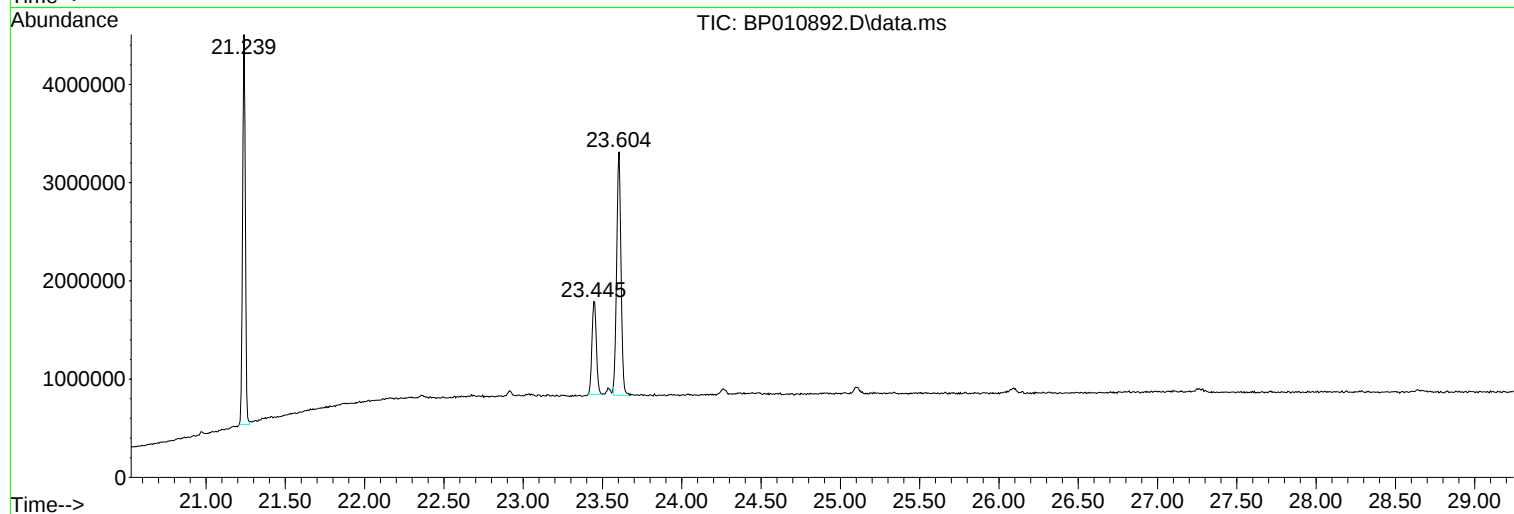
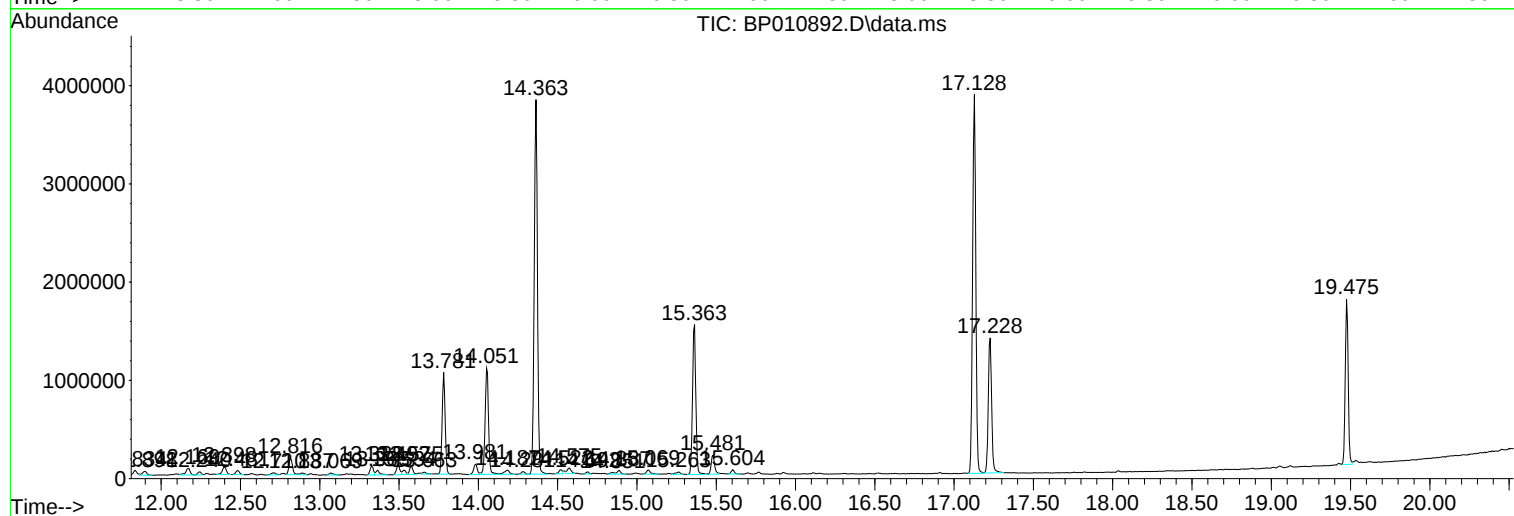
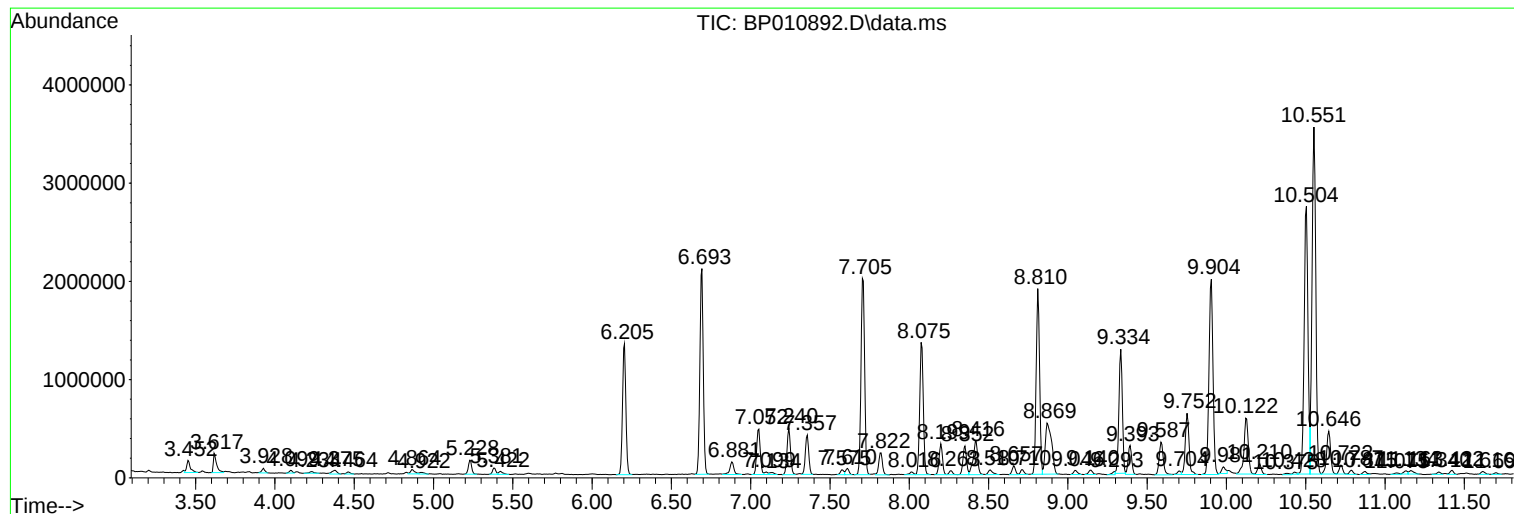
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Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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



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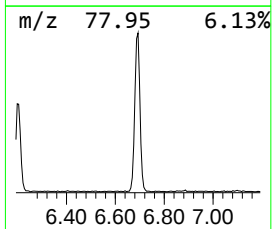
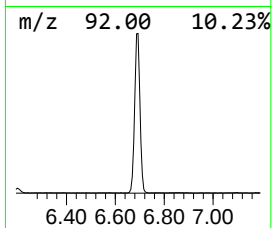
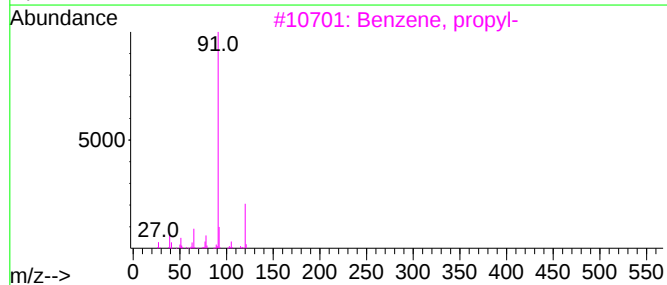
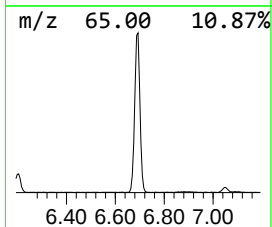
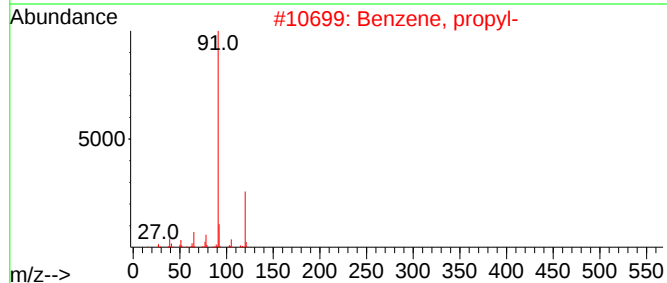
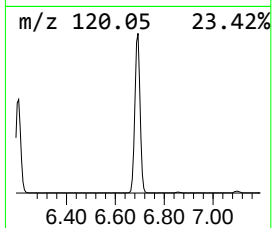
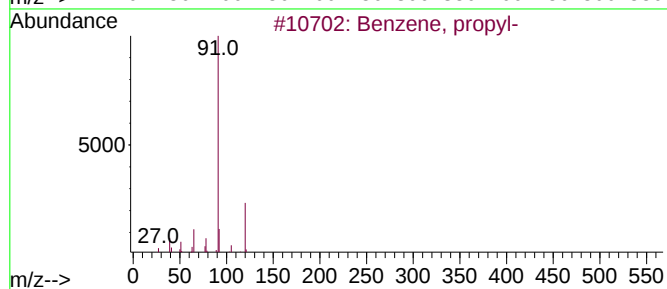
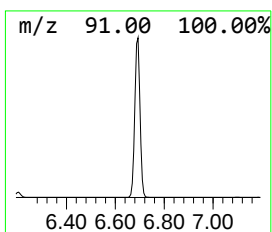
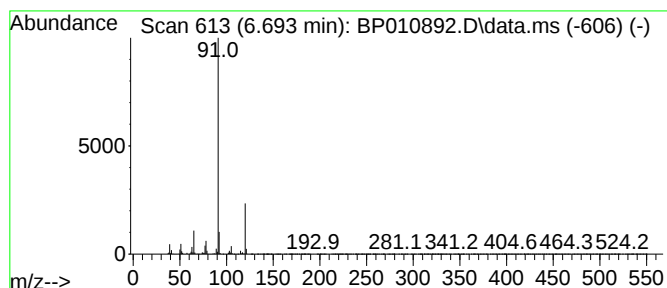
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 Benzene, propyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.693	19.50 ng/ul	3104820	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, propyl-	120	C9H12	000103-65-1	94
2			Benzene, propyl-	120	C9H12	000103-65-1	90
3			Benzene, propyl-	120	C9H12	000103-65-1	90
4			N-Benzyl-2-phenethylamine	211	C15H17N	003647-71-0	78
5			N-Isobutyl(phenyl)methanesulfona...	227	C11H17NO2S	144615-94-1	72



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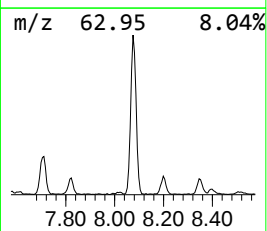
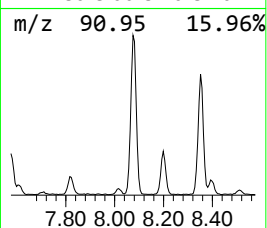
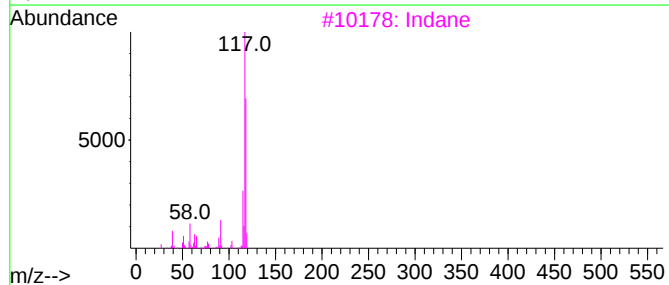
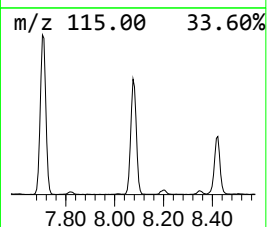
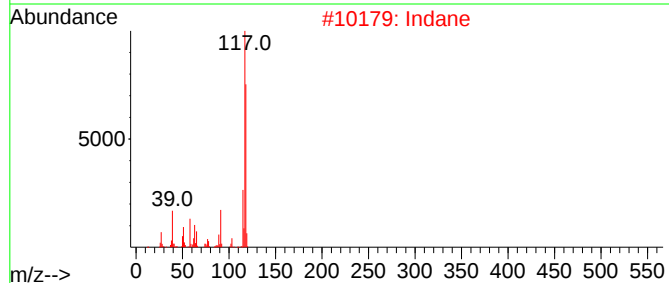
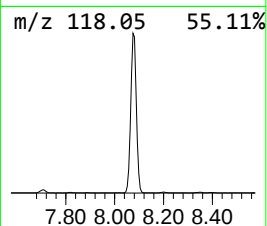
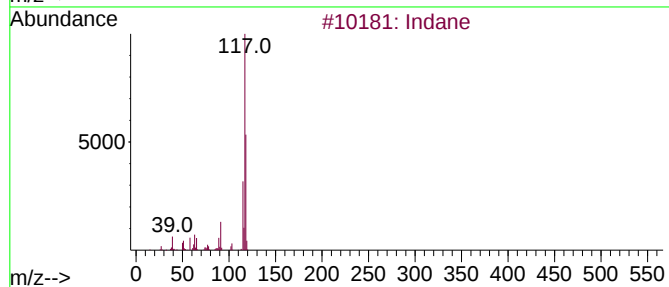
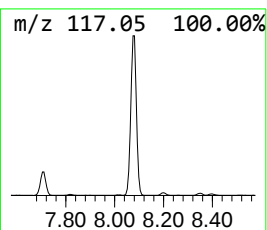
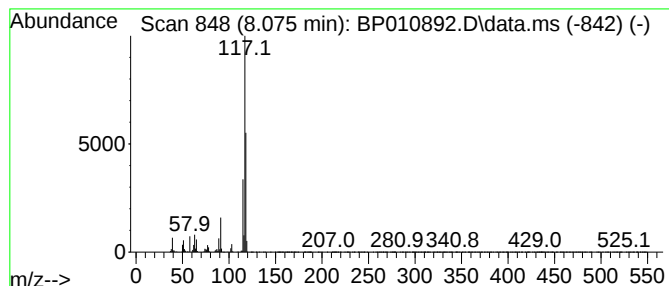
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 5 Indane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.075	13.33 ng/ul	2121820	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane			118	C9H10	000496-11-7	93
2	Indane			118	C9H10	000496-11-7	91
3	Indane			118	C9H10	000496-11-7	83
4	Deltacyclene			118	C9H10	007785-10-6	72
5	Indane			118	C9H10	000496-11-7	68



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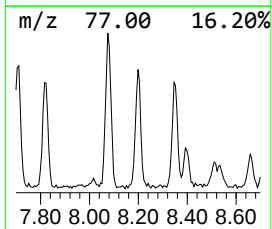
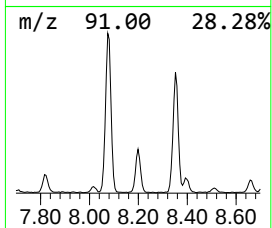
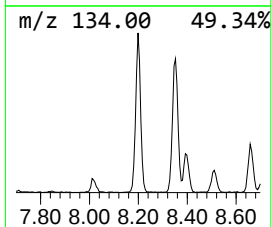
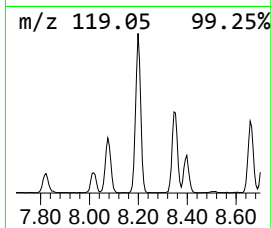
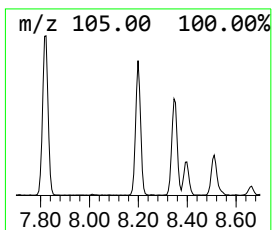
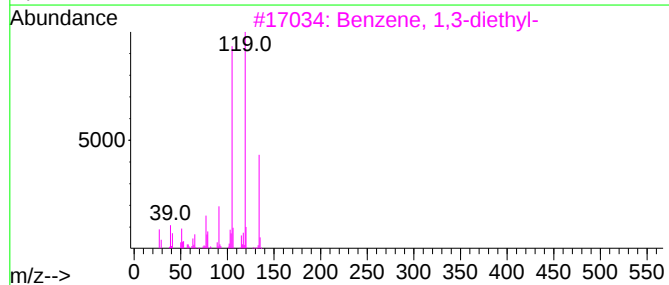
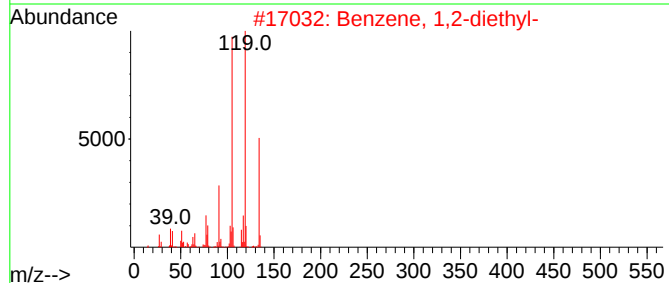
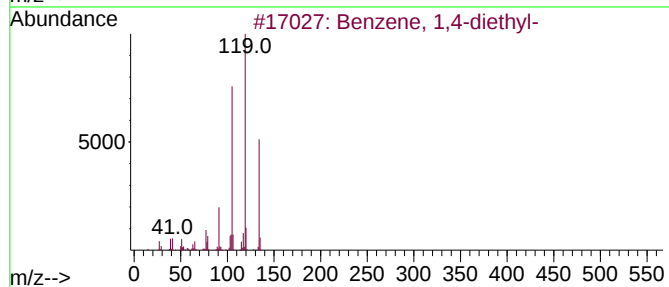
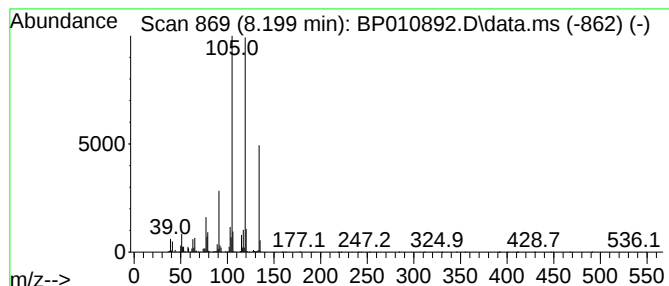
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.M
Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, 1,4-diethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.199	2.84 ng/ul	451750	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	97
2			Benzene, 1,2-diethyl-	134	C10H14	000135-01-3	96
3			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
4			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	96
5			Benzene, 1,4-diethyl-	134	C10H14	000105-05-5	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
Sample : N3401-14DL 5X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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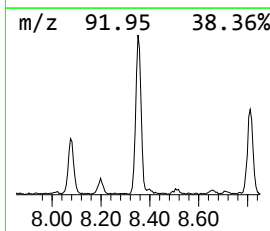
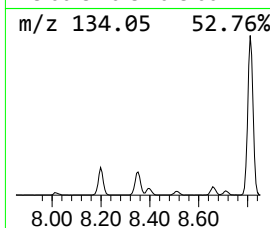
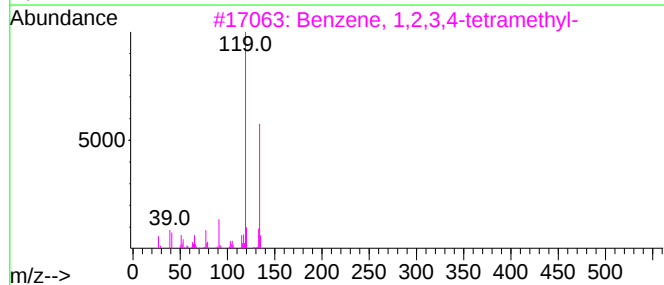
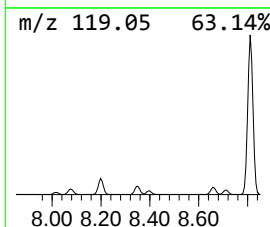
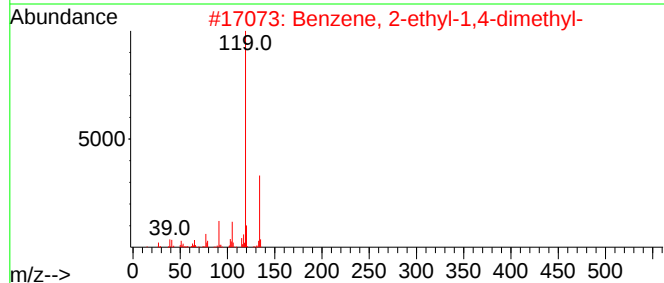
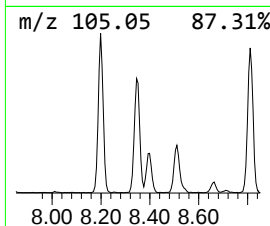
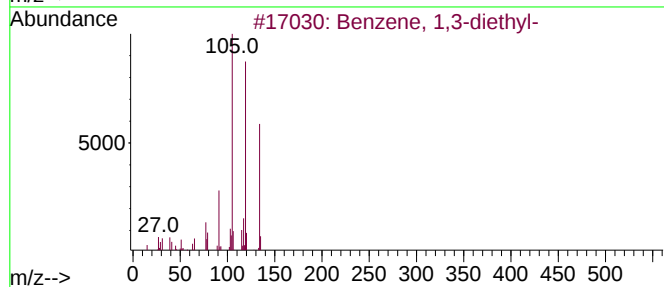
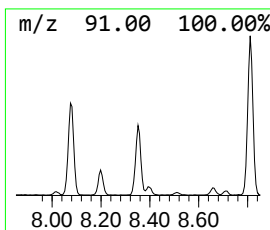
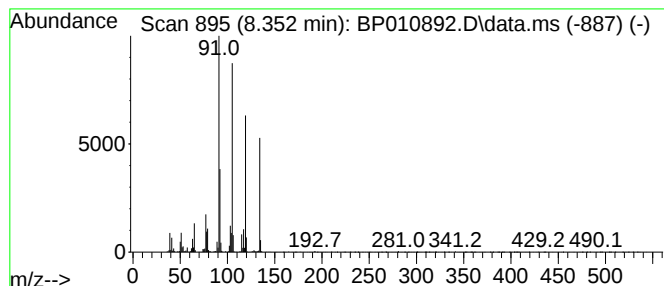
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, 1,3-diethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.352	2.87 ng/ul	456907	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,3-diethyl-	134	C10H14	000141-93-5	95
2			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	81
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	81
4			Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	76
5			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	76



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
Sample : N3401-14DL 5X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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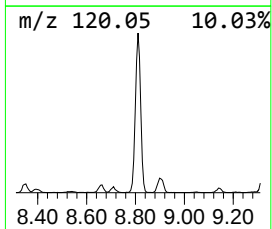
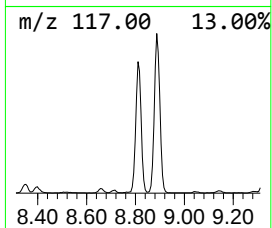
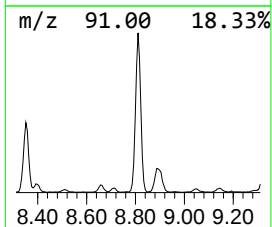
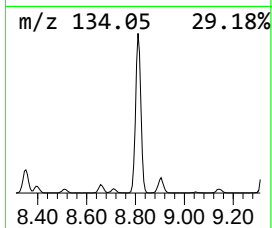
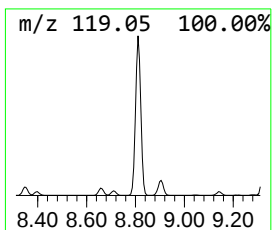
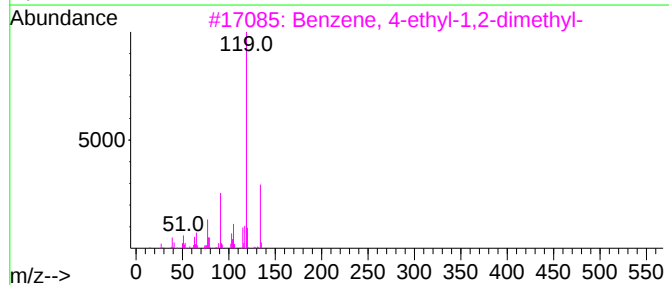
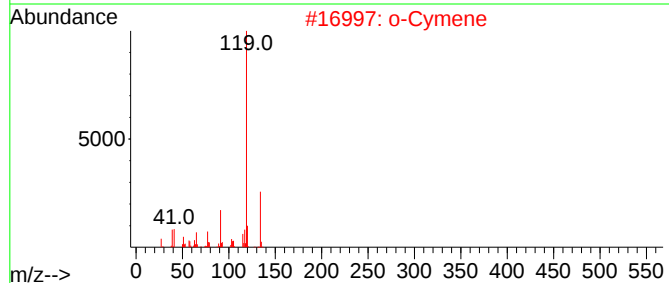
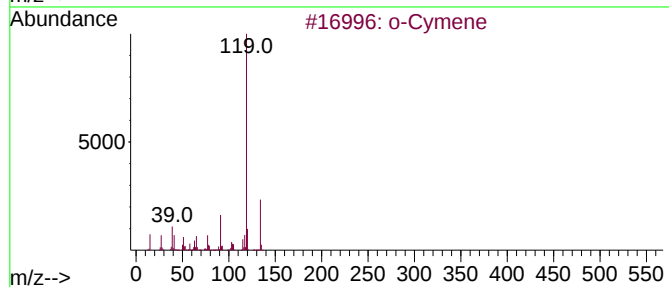
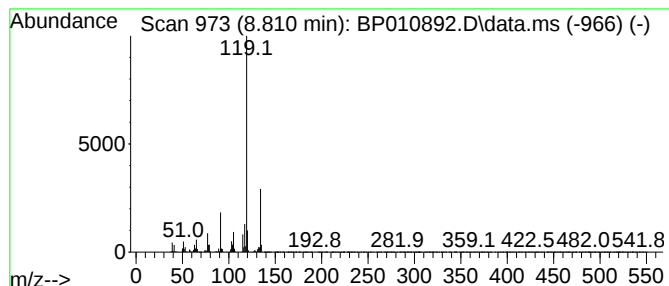
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 o-Cymene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.810	17.75 ng/ul	2825380	1,4-Dichlorobenzene-d4	7.705

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			o-Cymene	134	C10H14	000527-84-4	95
2			o-Cymene	134	C10H14	000527-84-4	95
3			Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
4			Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94
5			Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
Sample : N3401-14DL 5X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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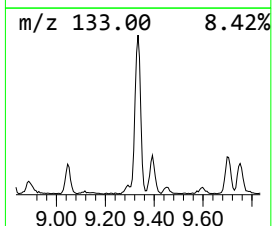
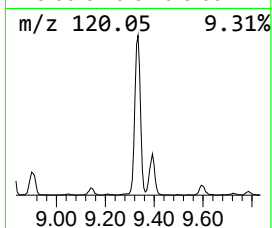
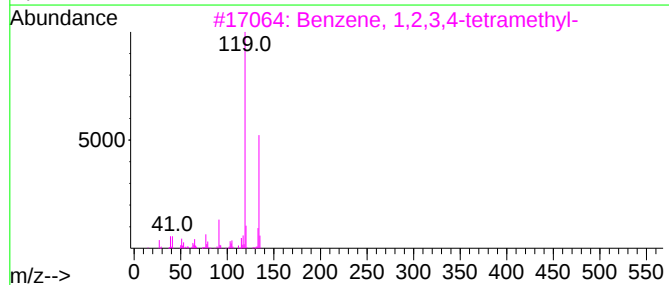
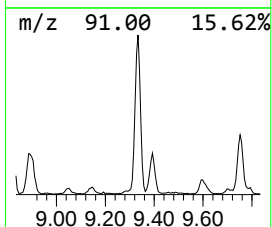
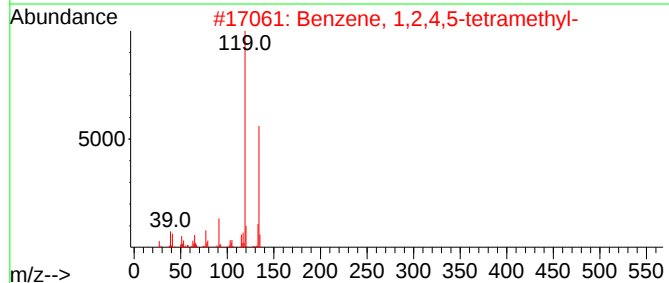
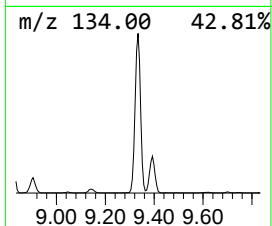
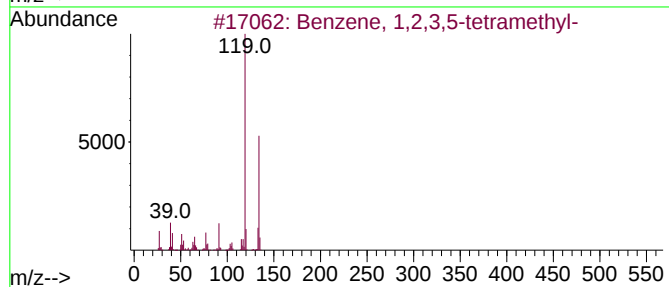
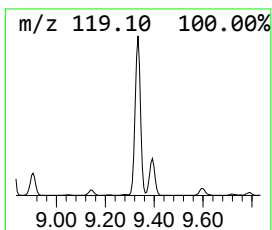
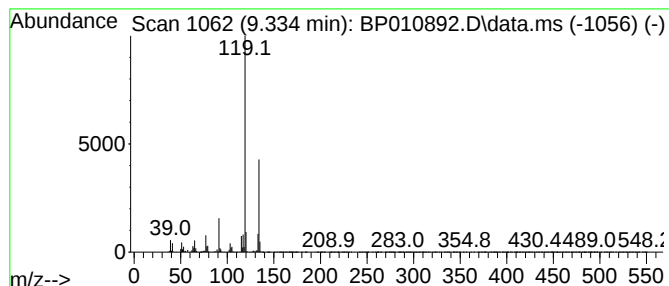
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, 1,2,3,5-tetramethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.334	8.15 ng/ul	1906130	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
2			Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	95
3			Benzene, 1,2,3,4-tetramethyl-	134	C10H14	000488-23-3	95
4			Benzene, 1,2,3,5-tetramethyl-	134	C10H14	000527-53-7	95
5			o-Cymene	134	C10H14	000527-84-4	95



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
Sample : N3401-14DL 5X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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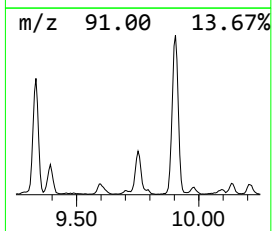
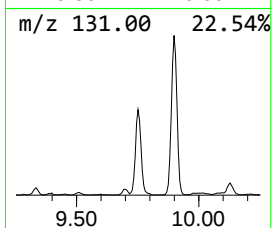
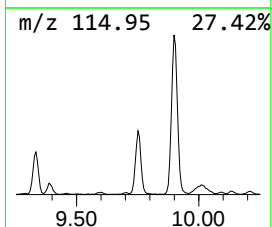
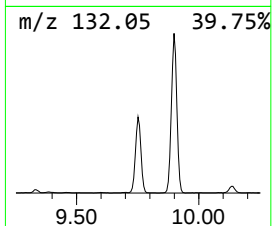
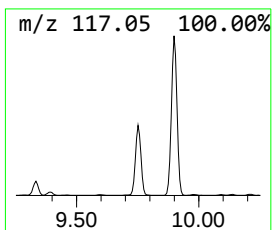
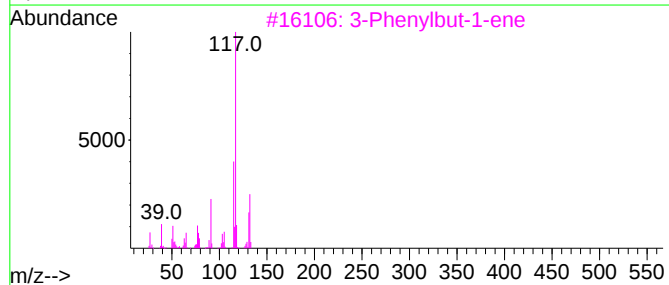
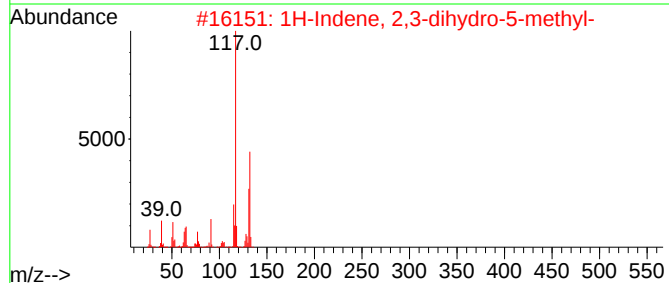
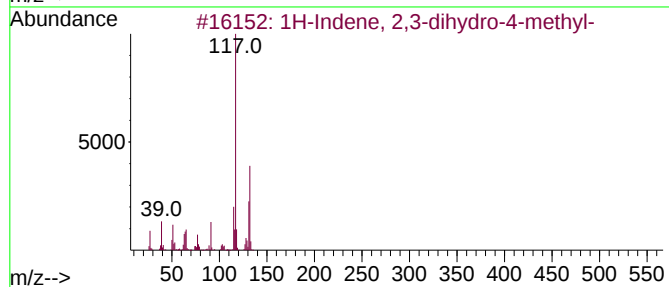
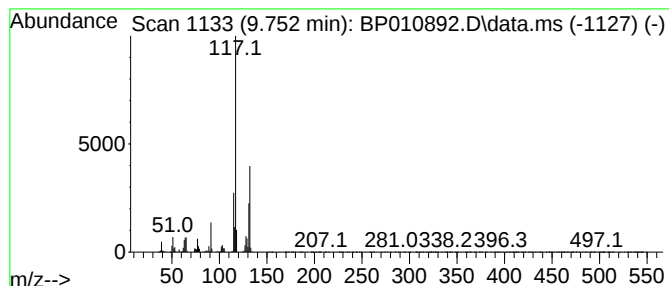
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 1H-Indene, 2,3-dihydro-4-me... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.752	4.32 ng/ul	1009690	Naphthalene-d8	10.504

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		1H-Indene, 2,3-dihydro-4-methyl-	132	C10H12	000824-22-6	94
2		1H-Indene, 2,3-dihydro-5-methyl-	132	C10H12	000874-35-1	93
3		3-Phenylbut-1-ene	132	C10H12	000934-10-1	90
4		Indan, 1-methyl-	132	C10H12	000767-58-8	87
5		Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	83



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
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Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
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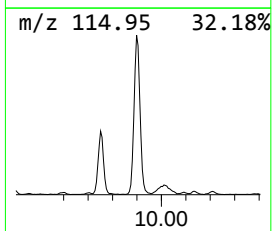
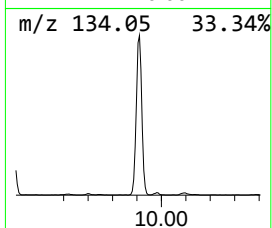
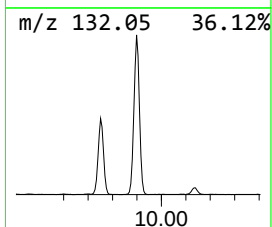
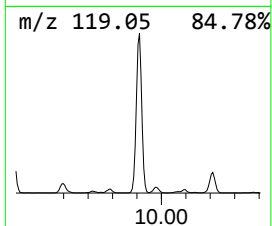
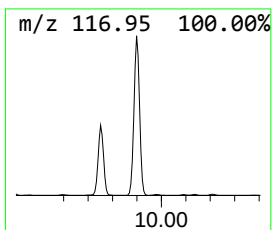
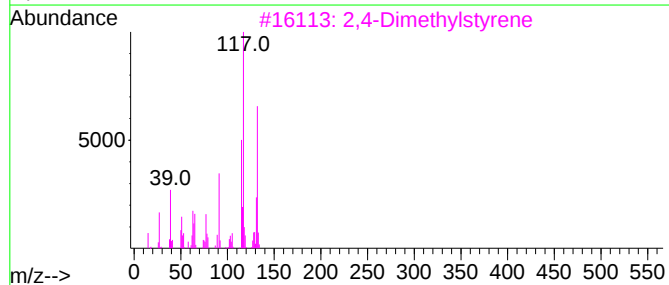
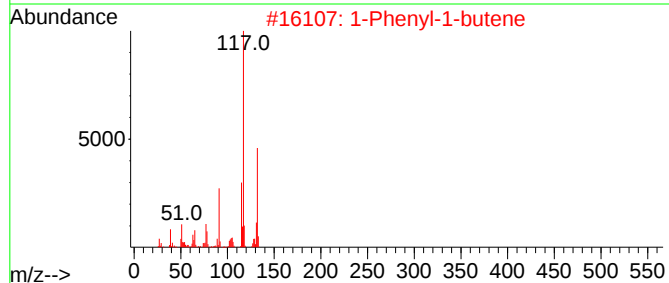
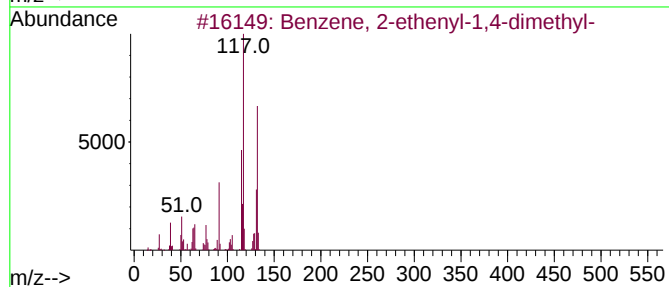
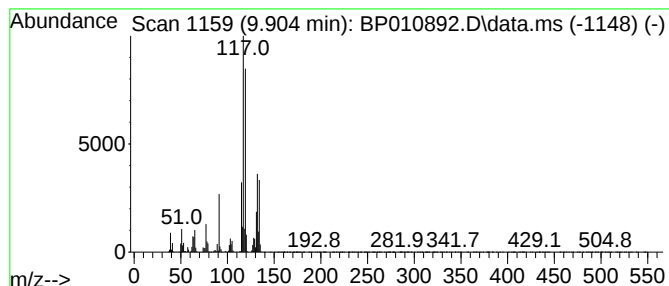
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.904	14.94 ng/ul	3493370	Naphthalene-d8	10.504

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	89
2			1-Phenyl-1-butene	132	C10H12	000824-90-8	87
3			2,4-Dimethylstyrene	132	C10H12	002234-20-0	86
4			3-Phenylbut-1-ene	132	C10H12	000934-10-1	70
5			Benzene, 2-butenyl-	132	C10H12	001560-06-1	55



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP062722\
Data File : BP010892.D
Acq On : 27 Jun 2022 23:53
Operator : CG/JU
Sample : N3401-14DL 5X
Misc :
ALS Vial : 25 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP062322.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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Indane	8.075	13.3	ng/u1	2121820	1	7.705	3184320	20.0
Benzene, 1,4-di...	8.199	2.8	ng/u1	451750	1	7.705	3184320	20.0
Benzene, 1,3-di...	8.352	2.9	ng/u1	456907	1	7.705	3184320	20.0
o-Cymene	8.810	17.8	ng/u1	2825380	1	7.705	3184320	20.0
Benzene, 1,2,3,...	9.334	8.2	ng/u1	1906130	2	10.504	4676420	20.0
1H-Indene, 2,3-...	9.752	4.3	ng/u1	1009690	2	10.504	4676420	20.0
Benzene, 2-ethe...	9.904	14.9	ng/u1	3493370	2	10.504	4676420	20.0