

Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
 Data File : BP016677.D
 Acq On : 21 Aug 2023 16:13
 Operator : MA/JU
 Sample : 04049-05
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_P
 ClientSampleId :
 BH9T5

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-BP081423.MA.M
 Title : SVOA CALIBRATION

Signal : TIC: BP016677.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	3.346	6	10	14	rBV	84188	119565	1.18%	0.127%
2	3.387	14	17	35	rVB	62697	123462	1.22%	0.131%
3	3.828	88	92	113	rBV	590517	963389	9.53%	1.019%
4	4.028	122	126	133	rVB	43481	70735	0.70%	0.075%
5	4.675	229	236	251	rVB	474832	759661	7.51%	0.804%
6	5.158	313	318	324	rVB4	15596	26079	0.26%	0.028%
7	5.399	351	359	365	rBV3	23393	45306	0.45%	0.048%
8	5.775	416	423	427	rBV3	14304	25250	0.25%	0.027%
9	5.840	429	434	436	rVV3	24750	38162	0.38%	0.040%
10	5.870	436	439	445	rVV2	48276	86361	0.85%	0.091%
11	5.922	445	448	453	rVV2	21856	34235	0.34%	0.036%
12	6.058	466	471	482	rVB7	14531	40488	0.40%	0.043%
13	6.228	495	500	507	rBV5	55656	120475	1.19%	0.127%
14	6.293	507	511	518	rVV	135120	219895	2.17%	0.233%
15	6.464	535	540	546	rVV	64078	102081	1.01%	0.108%
16	6.634	565	569	572	rBV	36050	63951	0.63%	0.068%
17	6.687	572	578	590	rVB5	67144	161558	1.60%	0.171%
18	6.793	590	596	602	rVB3	51433	88335	0.87%	0.093%
19	6.905	611	615	620	rVB7	14626	22720	0.22%	0.024%
20	6.987	620	629	635	rBV3	115766	226669	2.24%	0.240%
21	7.034	635	637	640	rVV2	22086	27212	0.27%	0.029%
22	7.075	640	644	649	rVB2	51318	75698	0.75%	0.080%
23	7.128	649	653	658	rBV	99230	150994	1.49%	0.160%
24	7.187	658	663	669	rVV	405867	632052	6.25%	0.669%
25	7.269	675	677	685	rVB8	10748	23791	0.24%	0.025%
26	7.381	690	696	707	rVB	1500898	2356292	23.31%	2.493%
27	7.481	707	713	720	rBV5	44384	89413	0.88%	0.095%
28	7.564	720	727	733	rBV	1337365	2194219	21.70%	2.322%
29	7.611	733	735	742	rVB	180582	241873	2.39%	0.256%
30	7.746	750	758	763	rBV5	32449	72038	0.71%	0.076%
31	7.916	779	787	795	rVB2	74795	168814	1.67%	0.179%
32	8.046	802	809	819	rBV	879957	1482815	14.67%	1.569%
33	8.164	825	829	835	rVB3	16328	29313	0.29%	0.031%
34	8.287	845	850	858	rBV9	16364	36678	0.36%	0.039%
35	8.375	860	865	872	rBV10	10577	29755	0.29%	0.031%
36	8.746	920	928	937	rBV	1116419	1840776	18.21%	1.948%

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Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
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37	8.816	937	940	944	rVB3	16721	23967	0.24%	0.025%
38	8.893	950	953	960	rVB9	14511	26656	0.26%	0.028%
39	9.122	982	992	999	rVB3	77864	153510	1.52%	0.162%
40	9.240	1005	1012	1025	rVB	1696202	2882635	28.51%	3.050%
41	9.381	1025	1036	1041	rBV	404322	683590	6.76%	0.723%
42	9.416	1041	1042	1048	rVB	65452	76939	0.76%	0.081%
43	9.593	1062	1072	1084	rVB	77304	157035	1.55%	0.166%
44	9.710	1084	1092	1101	rBV	31448	77856	0.77%	0.082%
45	9.840	1107	1114	1119	rVB	98127	152996	1.51%	0.162%
46	9.958	1121	1134	1145	rBV	1309478	2206697	21.83%	2.335%
47	10.122	1156	1162	1168	rVB	161695	261697	2.59%	0.277%
48	10.222	1174	1179	1188	rBV2	44413	88133	0.87%	0.093%
49	10.416	1204	1212	1214	rBV7	9043	21274	0.21%	0.023%
50	10.481	1215	1223	1233	rBV	1881619	3175301	31.41%	3.360%
51	10.769	1270	1272	1278	rVB3	15556	24199	0.24%	0.026%
52	10.875	1278	1290	1301	rBV	1311730	2324128	22.99%	2.459%
53	11.034	1309	1317	1331	rBV	1709436	2953426	29.21%	3.125%
54	11.146	1331	1336	1342	rVV	51396	89473	0.88%	0.095%
55	11.210	1342	1347	1360	rVB9	17207	37670	0.37%	0.040%
56	11.422	1376	1383	1392	rBV9	10339	24249	0.24%	0.026%
57	11.752	1434	1439	1444	rBV7	10376	20968	0.21%	0.022%
58	11.810	1446	1449	1456	rVB6	15021	27223	0.27%	0.029%
59	11.987	1470	1479	1483	rBV	38609	70739	0.70%	0.075%
60	12.022	1483	1485	1492	rVB6	18537	26793	0.27%	0.028%
61	12.140	1501	1505	1509	rBV	30438	44871	0.44%	0.047%
62	12.234	1515	1521	1528	rBV8	19203	42677	0.42%	0.045%
63	12.428	1549	1554	1556	rBV2	38805	58656	0.58%	0.062%
64	12.528	1563	1571	1577	rBV3	335575	638223	6.31%	0.675%
65	12.663	1589	1594	1600	rVB5	45798	74597	0.74%	0.079%
66	12.757	1605	1610	1626	rVB2	119378	244477	2.42%	0.259%
67	12.981	1638	1648	1657	rBV5	82926	248555	2.46%	0.263%
68	13.057	1657	1661	1666	rBV4	47521	76665	0.76%	0.081%
69	13.104	1666	1669	1674	rVB7	15353	23526	0.23%	0.025%
70	13.151	1674	1677	1684	rVB7	15427	32711	0.32%	0.035%
71	13.346	1701	1710	1716	rBV5	83896	228524	2.26%	0.242%
72	13.546	1740	1744	1748	rBV2	50257	78414	0.78%	0.083%
73	13.610	1749	1755	1760	rVB3	59137	133635	1.32%	0.141%
74	13.669	1760	1765	1772	rBV8	26754	54482	0.54%	0.058%
75	13.769	1779	1782	1786	rVB4	19414	25037	0.25%	0.026%
76	13.899	1801	1804	1809	rVB7	17588	28099	0.28%	0.030%

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77	14.110	1831	1840	1848	rBV	3776810	5204043	51.47%	5.506%
78	14.322	1873	1876	1881	rVB5	16438	23502	0.23%	0.025%
79	14.398	1881	1889	1901	rBV	4381595	6437748	63.67%	6.811%
80	14.504	1903	1907	1914	rVB2	42120	75268	0.74%	0.080%
81	14.698	1933	1940	1947	rBV	1990359	2888250	28.57%	3.056%
82	14.810	1955	1959	1963	rBV7	23393	36045	0.36%	0.038%
83	14.863	1963	1968	1969	rVV5	21802	30487	0.30%	0.032%
84	14.898	1969	1974	1985	rVB	243739	459691	4.55%	0.486%
85	15.075	1998	2004	2010	rBV3	104639	174340	1.72%	0.184%
86	15.234	2026	2031	2037	rBV2	99670	160688	1.59%	0.170%
87	15.693	2102	2109	2116	rBV	5790013	7922980	78.36%	8.383%
88	15.751	2116	2119	2123	rVV2	81122	126013	1.25%	0.133%
89	15.810	2123	2129	2145	rVV	1687662	2409768	23.83%	2.550%
90	15.928	2146	2149	2153	rVB3	26024	36441	0.36%	0.039%
91	17.463	2404	2410	2417	rBV	2581365	3508222	34.70%	3.712%
92	17.563	2421	2427	2439	rBV2	6095628	8789961	86.94%	9.300%
93	18.298	2548	2552	2565	rBV2	145103	285638	2.83%	0.302%
94	19.369	2731	2734	2739	rVB3	76883	95906	0.95%	0.101%
95	19.439	2742	2746	2750	rVB	69507	94056	0.93%	0.100%
96	19.534	2759	2762	2775	rBV3	124209	234902	2.32%	0.249%
97	19.798	2801	2807	2817	rBV	7803314	10110543	100.00%	10.697%
98	21.563	3102	3107	3113	rBV	2438192	3346998	33.10%	3.541%
99	23.974	3508	3517	3531	rVB2	3839751	8579708	84.86%	9.077%
100	24.139	3539	3545	3557	rVB	1447049	3075333	30.42%	3.254%

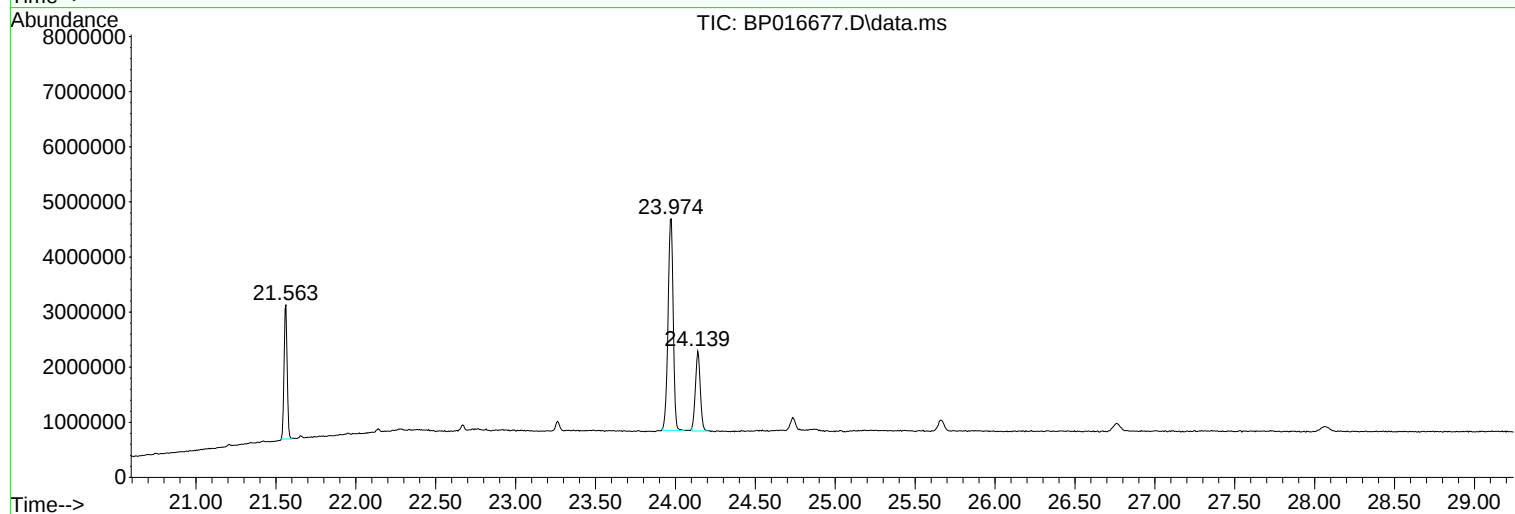
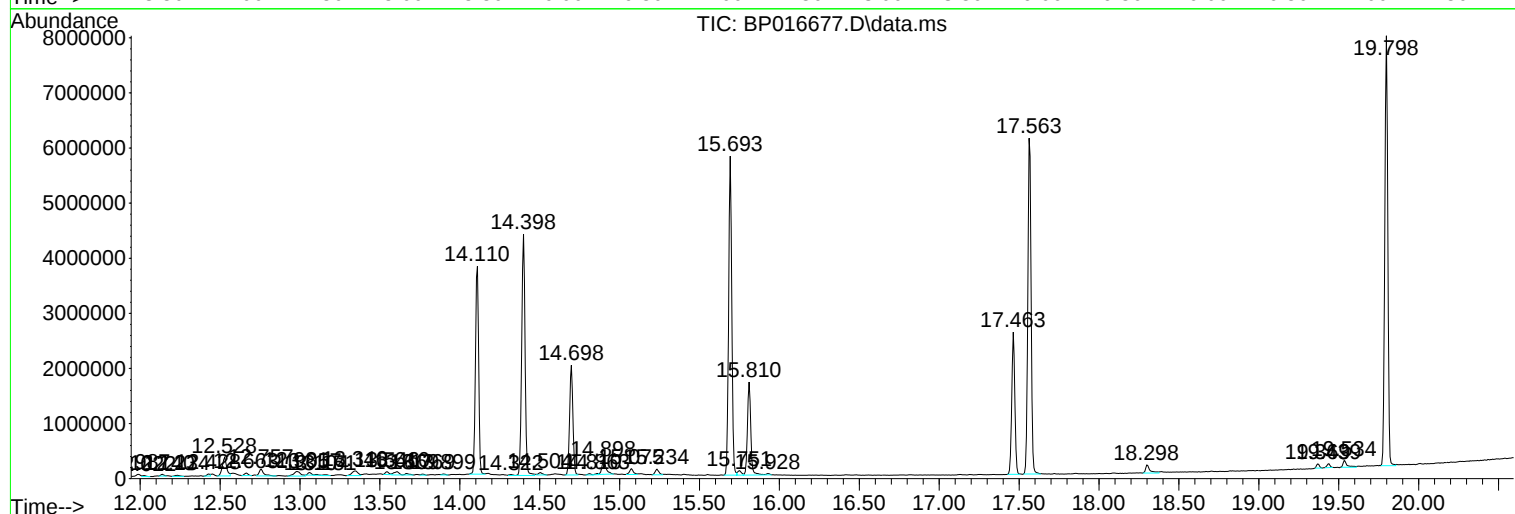
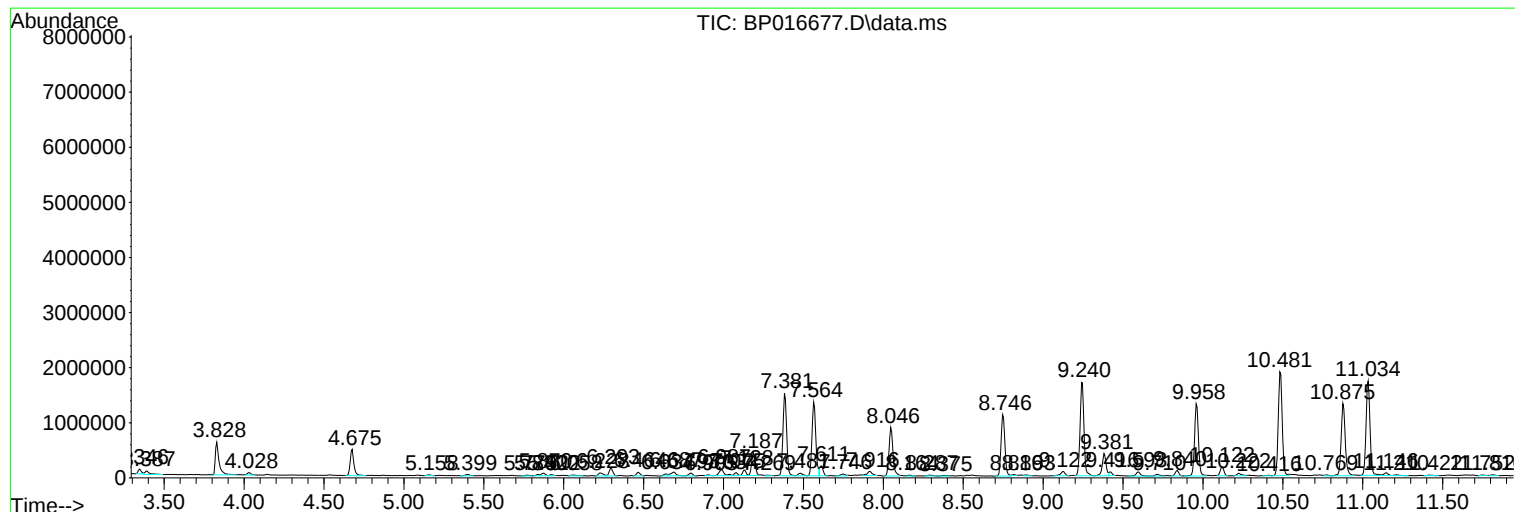
Sum of corrected areas: 94516944

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Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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



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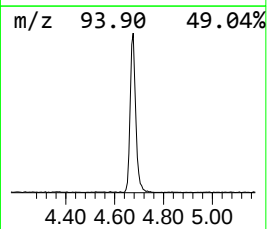
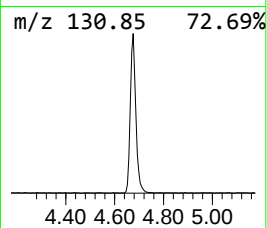
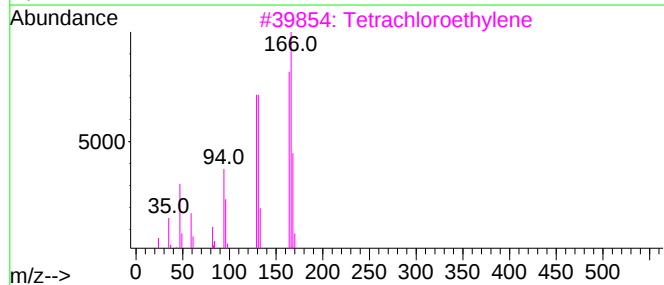
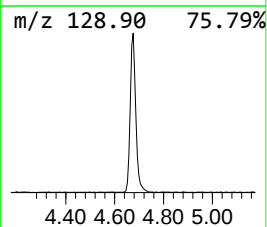
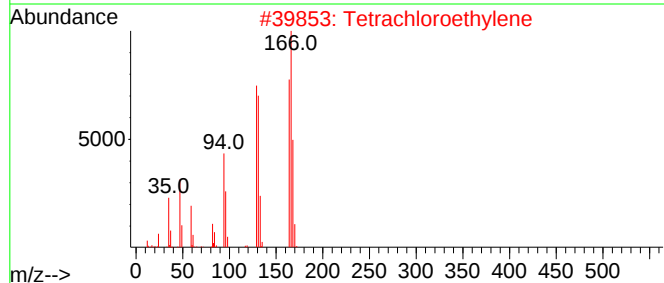
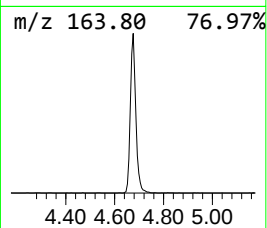
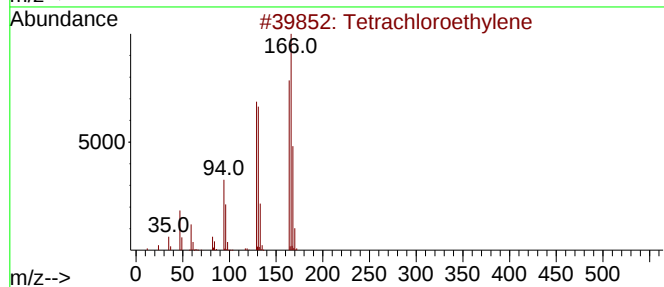
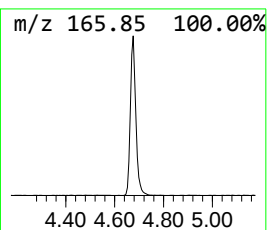
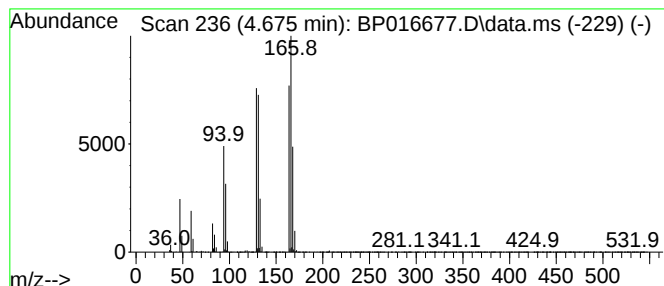
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 1 Tetrachloroethylene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
4.675	10.25 ng/ul	759661	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
2			Tetrachloroethylene	164	C2Cl4	000127-18-4	97
3			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
4			Tetrachloroethylene	164	C2Cl4	000127-18-4	96
5			Pyrimidine, 5-fluoro-2,4-dichloro-	166	C4HC12FN2	002927-71-1	12



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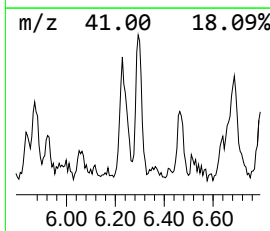
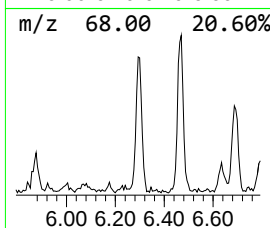
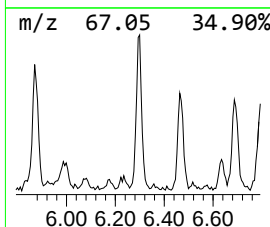
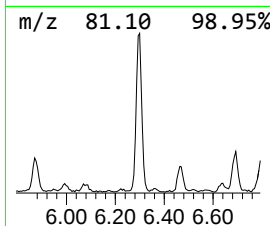
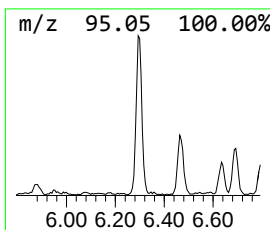
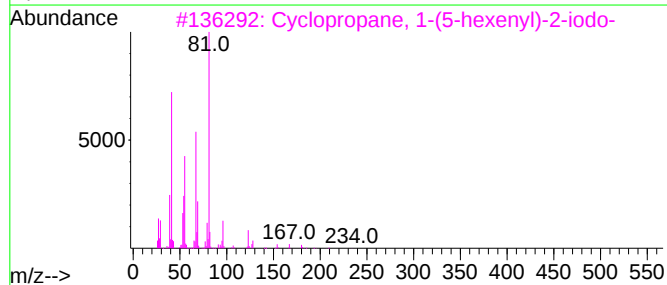
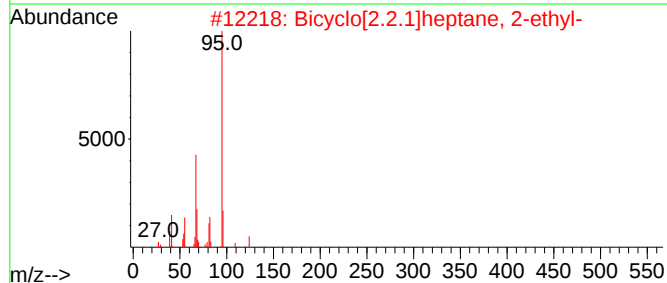
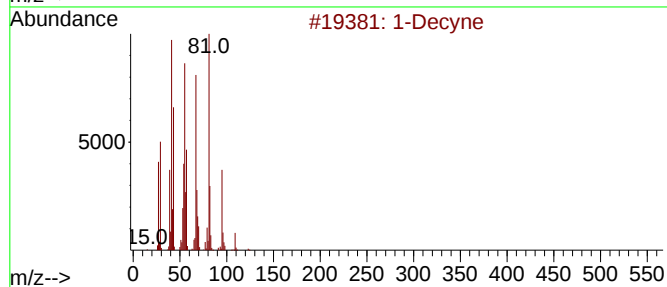
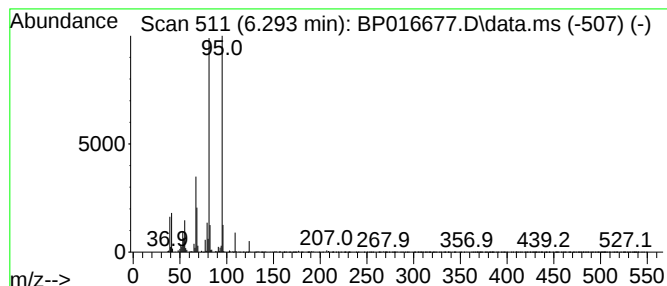
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 2 1-Decyne Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.293	2.97 ng/ul	219895	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Decyne			138	C10H18	000764-93-2	53
2	Bicyclo[2.2.1]heptane, 2-ethyl-			124	C9H16	002146-41-0	46
3	Cyclopropane, 1-(5-hexenyl)-2-iodo-			250	C9H15I	074685-40-8	43
4	Furan, 4-methyl-2-propyl-			124	C8H12O	006148-37-4	38
5	4-Methyl-1-prop-1-ynyl-cyclohexanol			152	C10H16O	1000362-94-7	38



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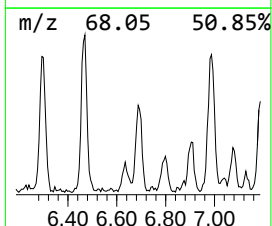
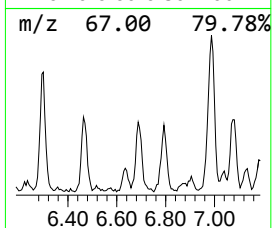
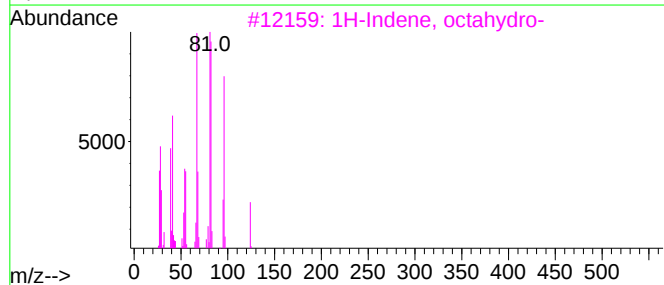
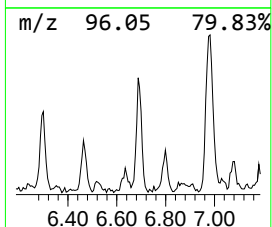
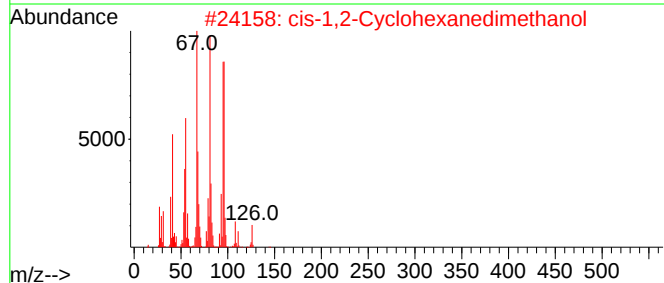
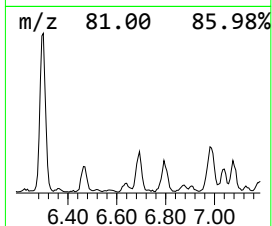
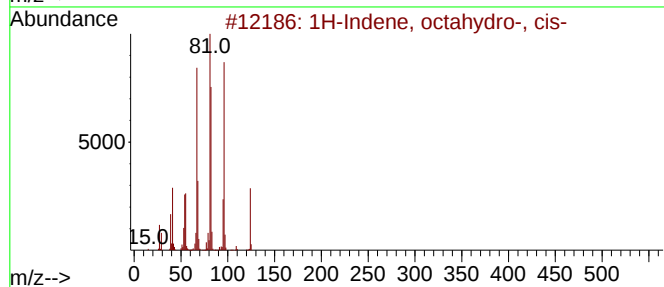
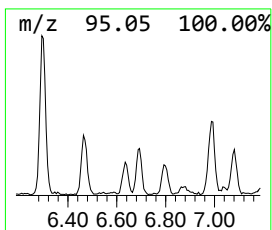
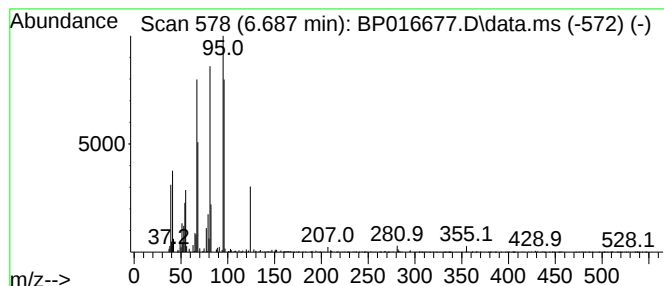
Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.MA.M
Quant Title : SVOA CALIBRATION

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

Peak Number 3 1H-Indene, octahydro-, cis- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.687	2.18 ng/ul	161558	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	68
2			cis-1,2-Cyclohexanedimethanol	144	C8H16O2	015753-50-1	64
3			1H-Indene, octahydro-	124	C9H16	000496-10-6	64
4			Cyclohexene, 1-octyl-	194	C14H26	015232-87-8	64
5			1H-Indene, octahydro-, cis-	124	C9H16	004551-51-3	62



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
Acq On : 21 Aug 2023 16:13
Operator : MA/JU
Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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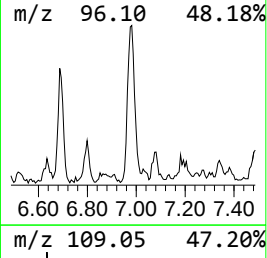
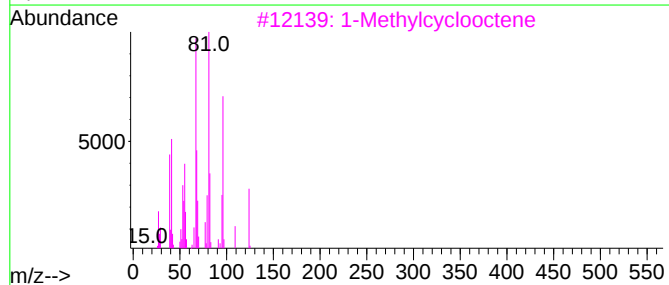
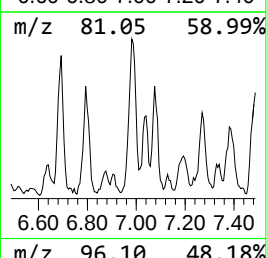
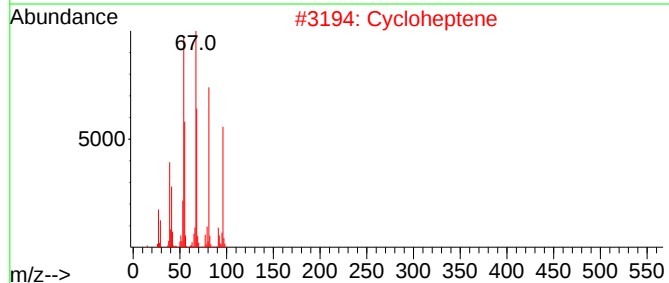
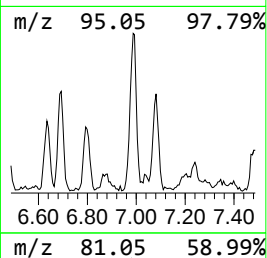
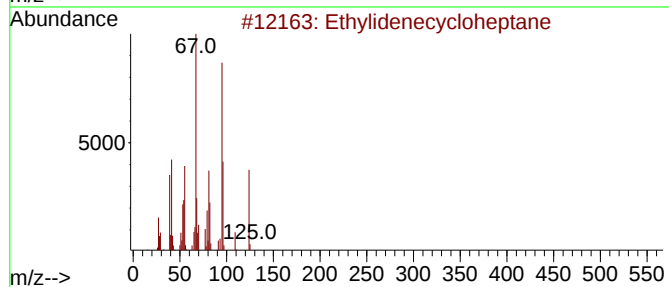
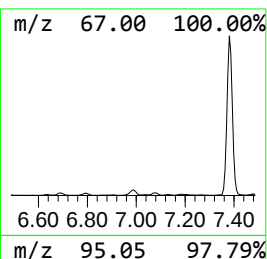
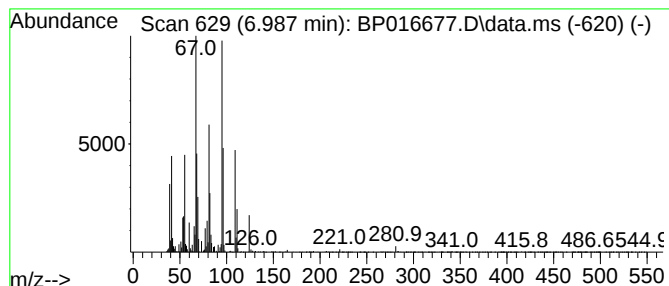
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Quant Title : SVOA CALIBRATION

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

Peak Number 4 Ethylenecycloheptane Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.987	3.06 ng/ul	226669	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Ethylenecycloheptane	124	C9H16	010494-87-8	86
2			Cycloheptene	96	C7H12	000628-92-2	83
3			1-Methylcyclooctene	124	C9H16	000933-11-9	83
4			2-Methylbicyclo[3.2.1]octane	124	C9H16	1010215-28-0	76
5			trans-1-Butenylcyclopentane	124	C9H16	1000113-50-9	60



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Data File : BP016677.D
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Operator : MA/JU
Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

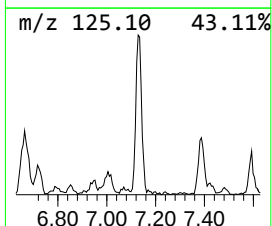
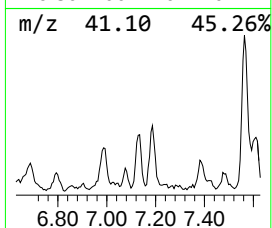
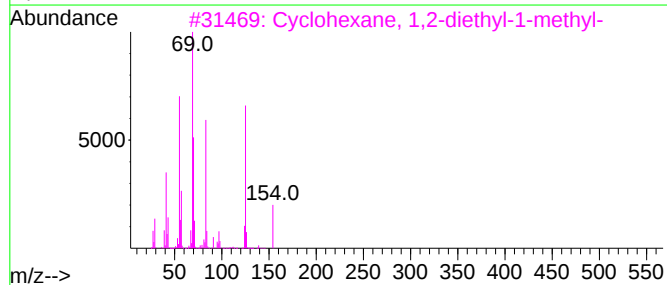
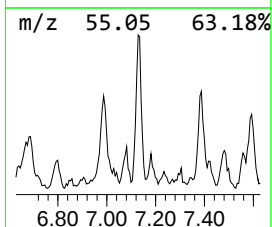
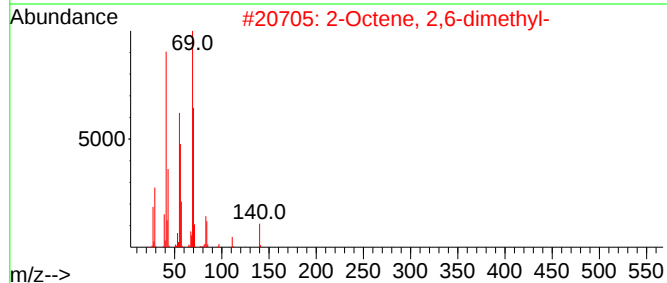
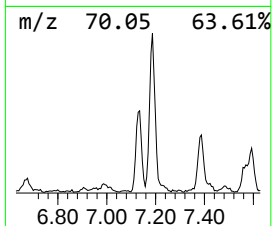
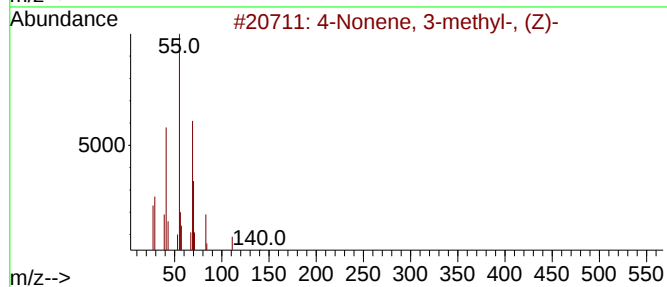
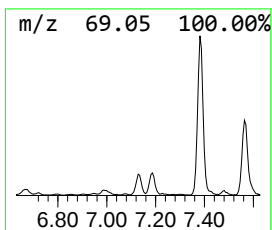
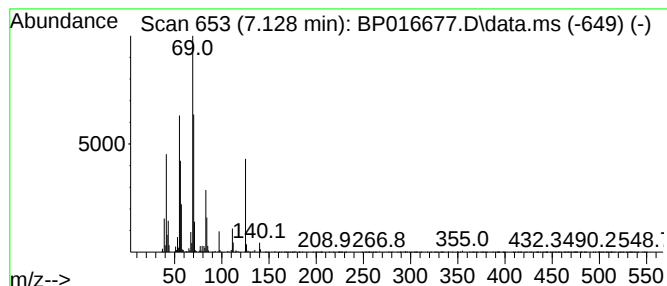
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Quant Title : SVOA CALIBRATION

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

Peak Number 5 (DEL) Alkane: Straight-Chai... Concentration Rank 11

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.128	2.04 ng/ul	150994	1,4-Dichlorobenzene-d4			8.046
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	4-Nonene, 3-methyl-, (Z)-		140	C10H20	063830-69-3	59
2	2-Octene, 2,6-dimethyl-		140	C10H20	004057-42-5	59
3	Cyclohexane, 1,2-diethyl-1-methyl-		154	C11H22	061141-79-5	53
4	Cyclopentane, ethyl-		98	C7H14	001640-89-7	47
5	2-n-Propylthiane, S,S-dioxide		176	C8H16O2S	072385-41-2	38



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Operator : MA/JU
Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

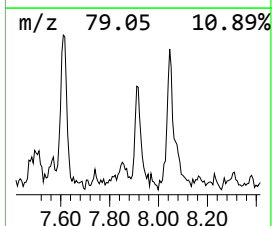
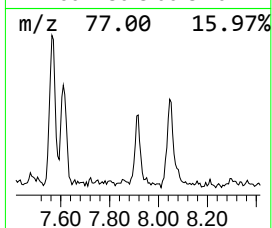
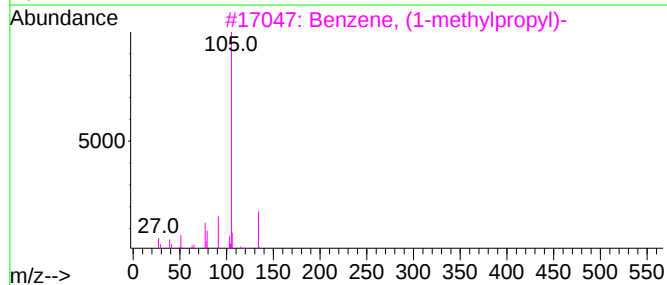
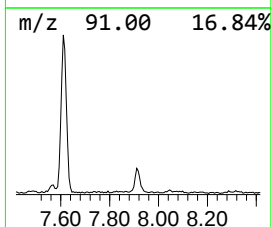
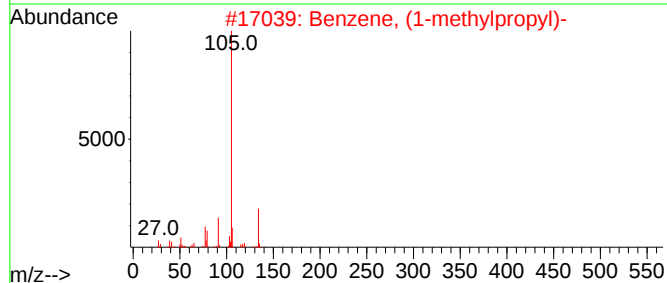
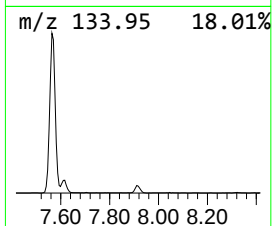
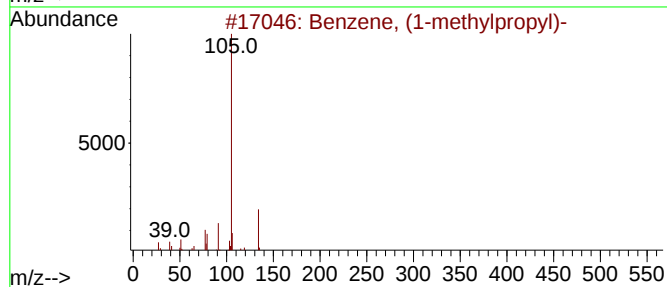
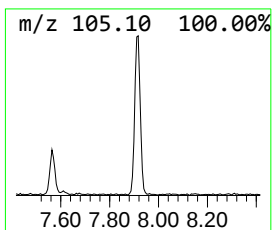
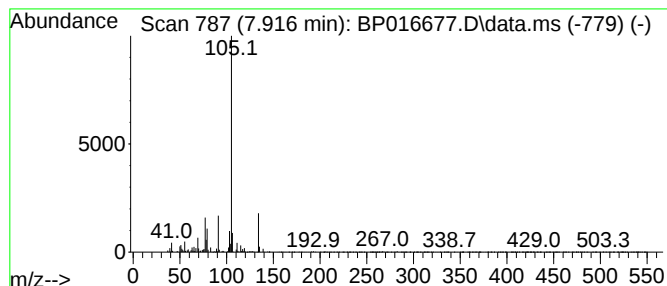
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Quant Title : SVOA CALIBRATION

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

Peak Number 6 Benzene, (1-methylpropyl)- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD			R.T.
7.916	2.28 ng/ul	168814	1,4-Dichlorobenzene-d4			8.046
Hit# of 5	Tentative ID		MW	MolForm	CAS#	Qual
1	Benzene, (1-methylpropyl)-		134	C10H14	000135-98-8	90
2	Benzene, (1-methylpropyl)-		134	C10H14	000135-98-8	90
3	Benzene, (1-methylpropyl)-		134	C10H14	000135-98-8	90
4	Benzene, (1-methylpropyl)-		134	C10H14	000135-98-8	90
5	Benzeneacetaldehyde, .alpha.-met...		134	C9H10O	000093-53-8	87



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
Acq On : 21 Aug 2023 16:13
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Misc :
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Instrument :
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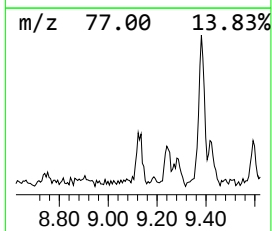
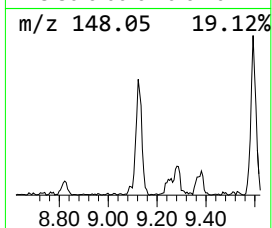
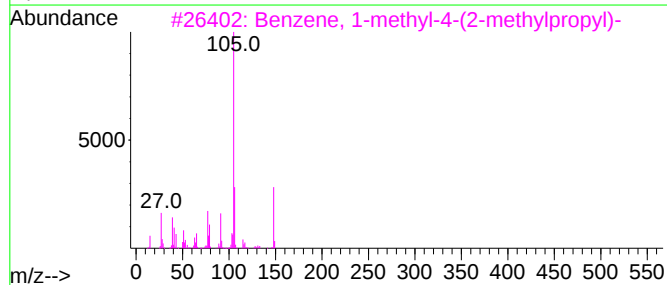
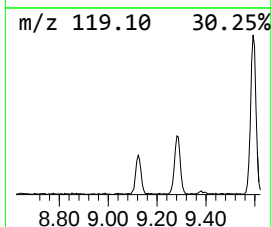
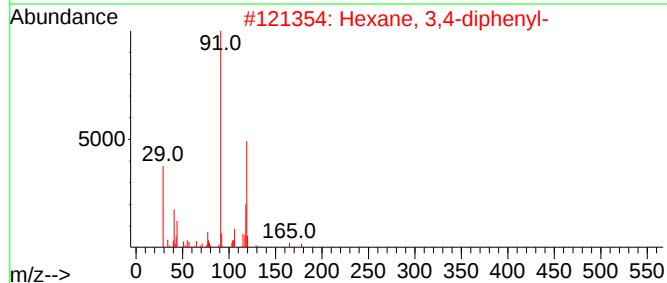
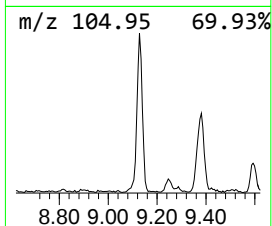
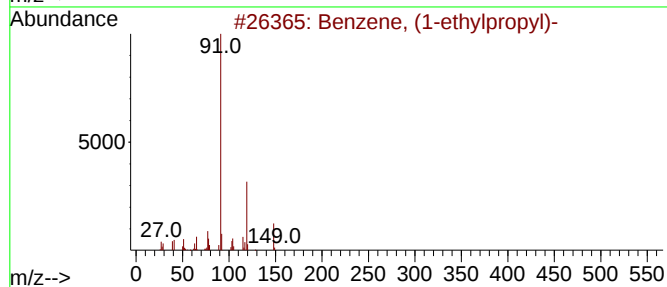
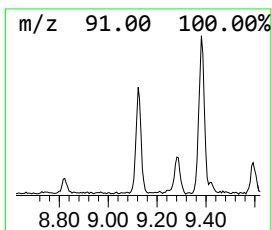
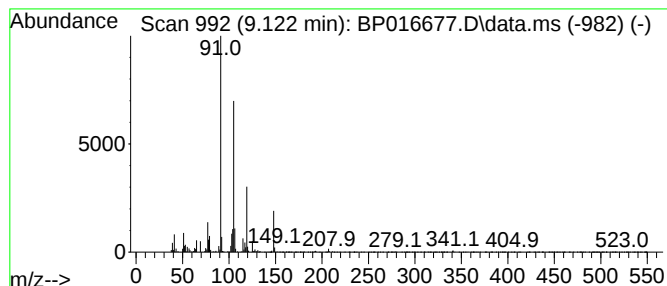
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Quant Title : SVOA CALIBRATION

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

Peak Number 7 Benzene, (1-ethylpropyl)- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.122	2.07 ng/ul	153510	1,4-Dichlorobenzene-d4	8.046

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-ethylpropyl)-	148	C11H16	001196-58-3	62
2			Hexane, 3,4-diphenyl-	238	C18H22	005789-31-1	43
3			Benzene, 1-methyl-4-(2-methylpro...	148	C11H16	005161-04-6	38
4			Benzene, (1-nitropropyl)-	165	C9H11NO2	005279-14-1	38
5			Benzene, (1-methylbutyl)-	148	C11H16	002719-52-0	35



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
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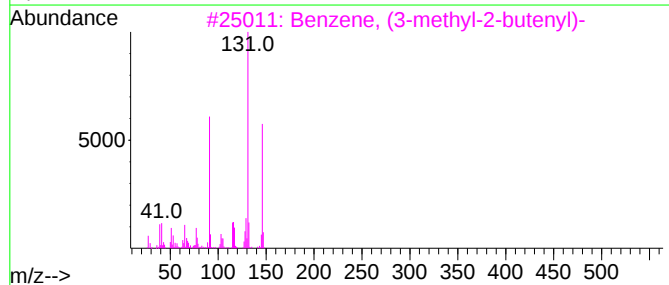
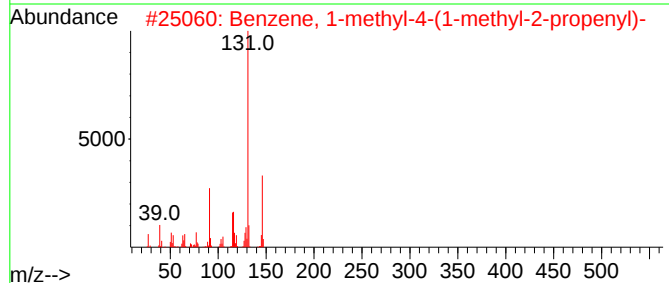
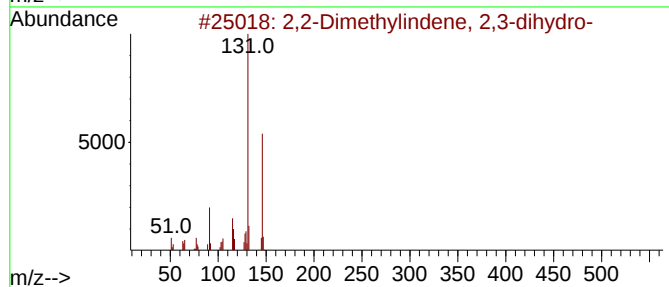
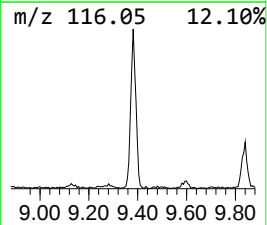
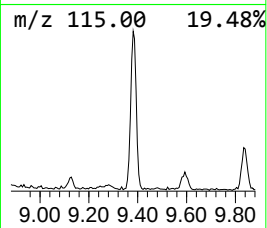
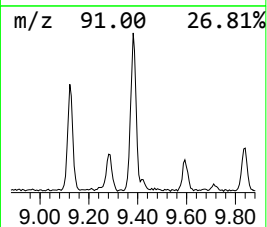
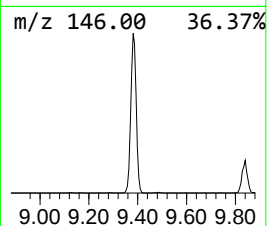
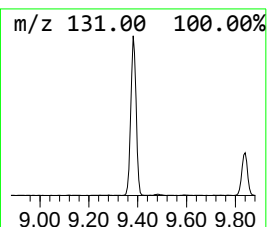
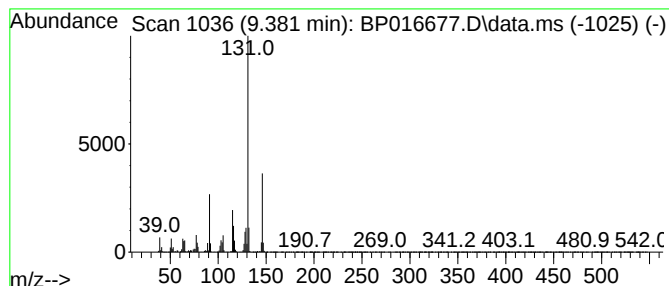
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Quant Title : SVOA CALIBRATION

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

Peak Number 8 2,2-Dimethylindene, 2,3-dih... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.381	9.22 ng/ul	683590	1,4-Dichlorobenzene-d4	8.046

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	97
2		Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	95
3		Benzene, (3-methyl-2-butenyl)-	146	C11H14	004489-84-3	94
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	94
5		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
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Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

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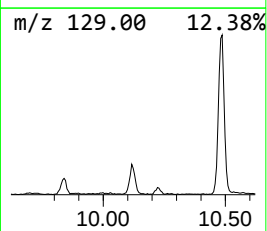
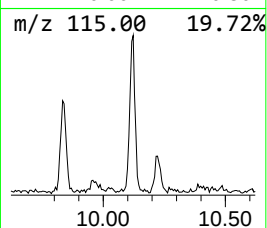
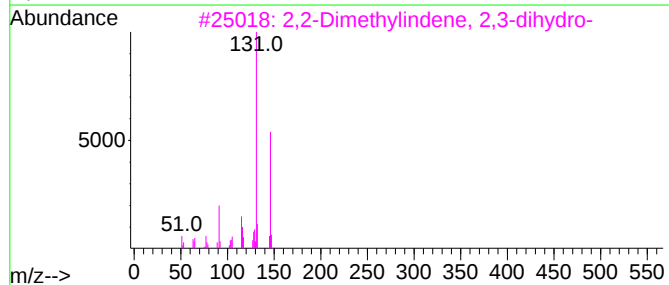
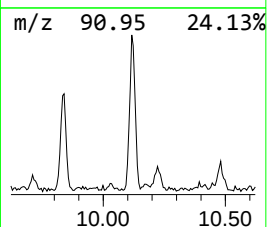
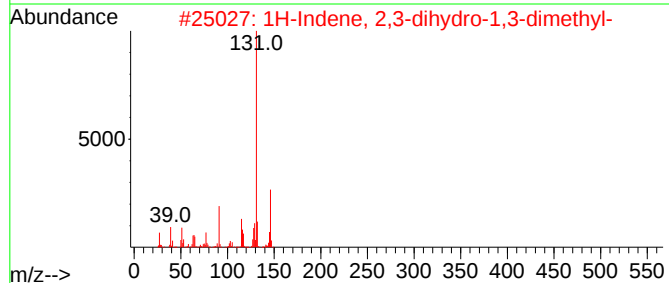
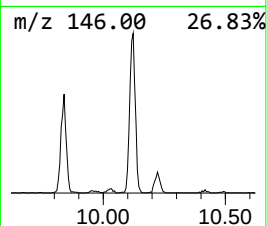
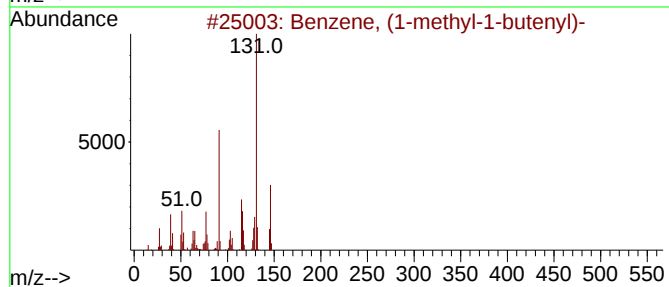
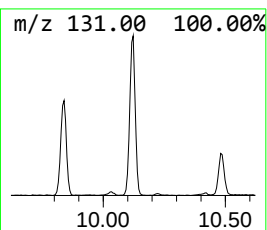
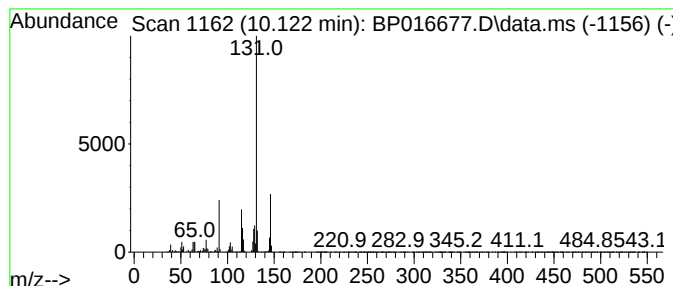
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Quant Title : SVOA CALIBRATION

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

Peak Number 9 Benzene, (1-methyl-1-butenyl)- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.122	2.25 ng/ul	261697	Naphthalene-d8	10.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	97
2			1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	94
3			2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	94
4			Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	94
5			1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
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Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
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ClientSampleId :
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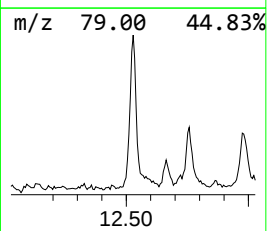
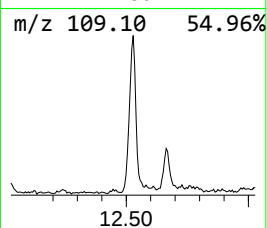
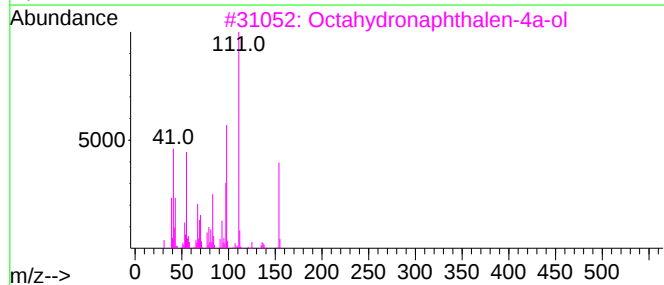
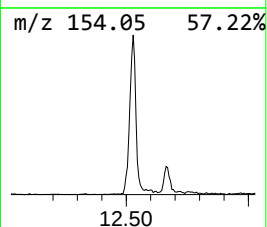
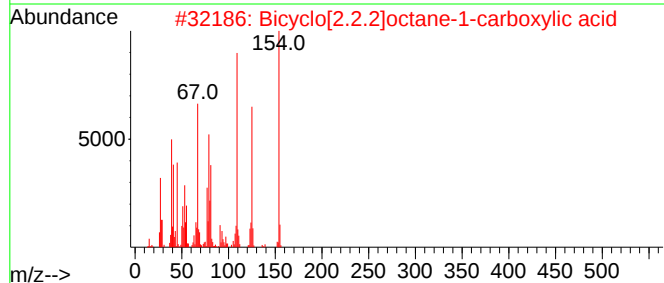
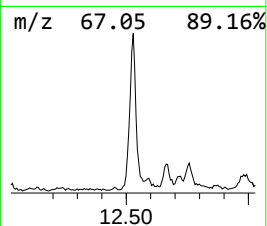
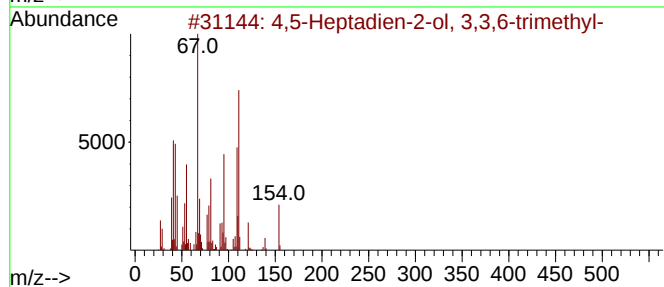
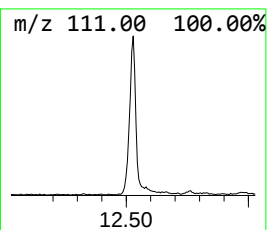
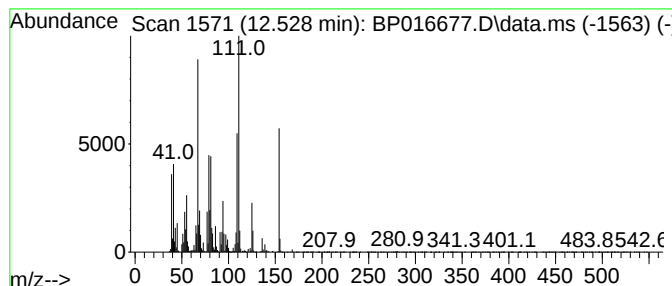
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Quant Title : SVOA CALIBRATION

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

Peak Number 10 4,5-Heptadien-2-ol, 3,3,6-t... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.528	5.49 ng/ul	638223	Naphthalene-d8	10.875

Hit#	of 5	Tentative ID	MW	MolForm	CAS#	Qual
1		4,5-Heptadien-2-ol, 3,3,6-trimet...	154	C10H18O	060432-69-1	58
2		Bicyclo[2.2.2]octane-1-carboxyli...	154	C9H14O2	000699-55-8	43
3		Octahydronaphthalen-4a-ol	154	C10H18O	055693-34-0	38
4		Cyclopentene-1-carboxamide	111	C6H9NO	1000423-09-9	35
5		2,4-Heptadiene, 2,4-dimethyl-	124	C9H16	074421-05-9	22



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
Acq On : 21 Aug 2023 16:13
Operator : MA/JU
Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH9T5

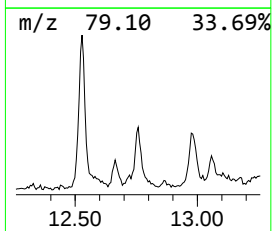
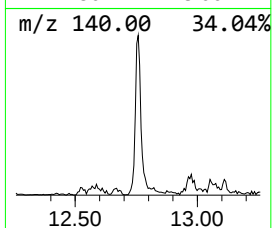
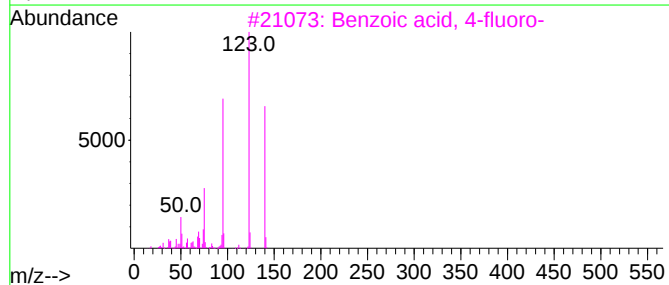
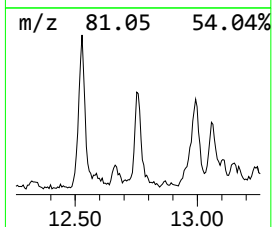
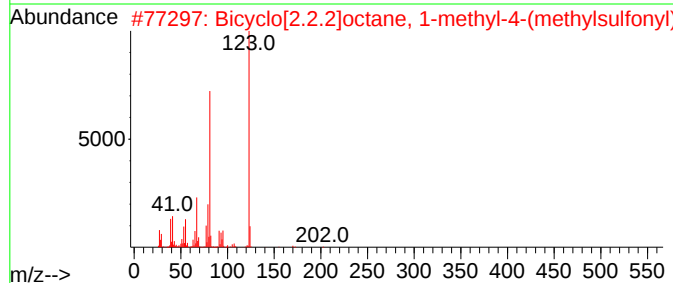
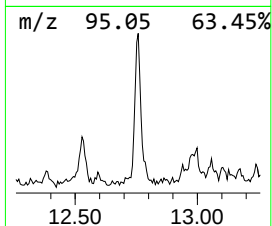
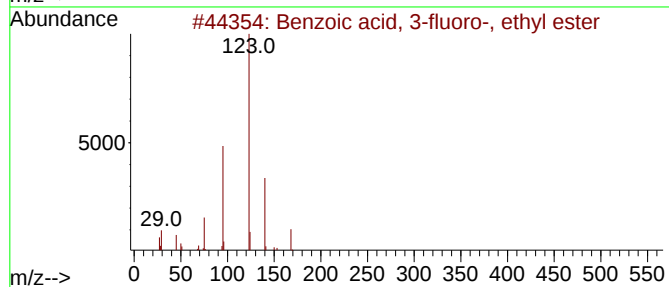
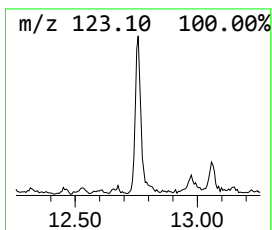
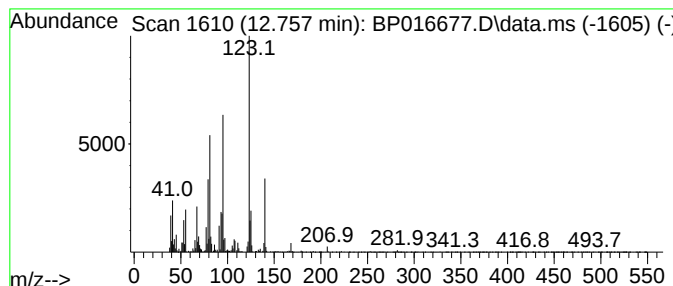
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Quant Title : SVOA CALIBRATION

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

Peak Number 11 Benzoic acid, 3-fluoro-, et... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.757	2.10 ng/ul	244477	Naphthalene-d8	10.875

Hit#	of	5	Tentative ID	MW	MolForm	CAS#	Qual
1			Benzoic acid, 3-fluoro-, ethyl e...	168	C9H9F02	000451-02-5	58
2			Bicyclo[2.2.2]octane, 1-methyl-4...	202	C10H18O2S	069855-48-7	55
3			Benzoic acid, 4-fluoro-	140	C7H5F02	000456-22-4	52
4			2,2-Dimethylocta-3,4-dienal	152	C10H16O	000590-71-6	52
5			Benzoic acid, 4-fluoro-	140	C7H5F02	000456-22-4	52



Data Path : Z:\svoasrv\HPCHEM1\BNA_P\Data\BP082123\
Data File : BP016677.D
Acq On : 21 Aug 2023 16:13
Operator : MA/JU
Sample : 04049-05
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument :
BNA_P
ClientSampleId :
BH9T5

Quant Method : Z:\svoasrv\HPCHEM1\BNA_P\Methods\SFAM-EPA-BP081423.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
Tetrachloroethy...	4.675	10.3	ng/ul	759661	1	8.046	1482820	20.0
1-Decyne	6.293	3.0	ng/ul	219895	1	8.046	1482820	20.0
1H-Indene, octa...	6.687	2.2	ng/ul	161558	1	8.046	1482820	20.0
Ethylidenecyclo...	6.987	3.1	ng/ul	226669	1	8.046	1482820	20.0
(DEL) Alkane: S...	7.128	2.0	ng/ul	150994	1	8.046	1482820	20.0
Benzene, (1-met...	7.916	2.3	ng/ul	168814	1	8.046	1482820	20.0
Benzene, (1-eth...	9.122	2.1	ng/ul	153510	1	8.046	1482820	20.0
2,2-Dimethylind...	9.381	9.2	ng/ul	683590	1	8.046	1482820	20.0
Benzene, (1-met...	10.122	2.3	ng/ul	261697	2	10.875	2324130	20.0
4,5-Heptadien-2...	12.528	5.5	ng/ul	638223	2	10.875	2324130	20.0
Benzoic acid, 3...	12.757	2.1	ng/ul	244477	2	10.875	2324130	20.0