

Data Path : Z:\VOASRV\HPCHEM1\MSVOA F\DATA\VF060719\
 Data File : VF062898.D
 Acq On : 7 Jun 2019 14:23
 Operator : FY/SY
 Sample : K3160-01
 Misc : 7.75g/5mL/MSVOA F/SOIL
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_F
 ClientSampleId :
 EO-02-060619-A

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
 Title : SW846 8260

Signal : TIC

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	1.065	52	54	56	rBV2	18352	35421	1.48%	0.079%
2	1.115	58	59	62	rVV	25276	46525	1.94%	0.104%
3	2.297	176	179	180	rBV3	24618	42803	1.78%	0.096%
4	2.514	197	201	204	rVB6	16203	36236	1.51%	0.081%
5	3.125	262	263	267	rVB	27177	41352	1.72%	0.093%
6	4.229	368	375	378	rBV4	14686	49576	2.07%	0.111%
7	4.318	378	384	391	rVV	363243	1095324	45.67%	2.455%
8	4.841	432	437	439	rBV4	17697	38126	1.59%	0.085%
9	5.067	452	460	469	rBV2	677631	2288936	95.44%	5.130%
10	5.787	526	533	550	rVB	898594	2398233	100.00%	5.375%
11	6.871	642	643	644	rBV	51143	35295	1.47%	0.079%
12	7.453	697	702	706	rBV8	14063	44071	1.84%	0.099%
13	7.719	722	729	743	rBV	838483	2021273	84.28%	4.530%
14	8.419	797	800	805	rVB4	15779	33601	1.40%	0.075%
15	8.764	831	835	838	rBV4	12063	25155	1.05%	0.056%
16	9.641	920	924	928	rVB3	18832	41979	1.75%	0.094%
17	9.917	944	952	962	rBV	1061323	2305657	96.14%	5.167%
18	10.035	962	964	969	rVB4	10899	30075	1.25%	0.067%
19	10.262	983	987	991	rBV4	10564	25303	1.06%	0.057%
20	10.824	1038	1044	1048	rVB2	28571	58177	2.43%	0.130%
21	10.893	1048	1051	1056	rBV6	10454	29301	1.22%	0.066%
22	11.455	1104	1108	1109	rBV2	26369	43567	1.82%	0.098%
23	11.504	1109	1113	1119	rVV	618897	1100382	45.88%	2.466%
24	11.632	1123	1126	1128	rVV4	21068	37717	1.57%	0.085%
25	11.682	1128	1131	1134	rVB	21446	39011	1.63%	0.087%
26	11.800	1139	1143	1150	rVV	66062	158706	6.62%	0.356%
27	11.898	1150	1153	1157	rVB	84890	128820	5.37%	0.289%
28	11.977	1157	1161	1165	rBV6	13179	46405	1.93%	0.104%
29	12.056	1167	1169	1170	rVV2	22209	32944	1.37%	0.074%
30	12.095	1170	1173	1176	rVB	94601	147211	6.14%	0.330%
31	12.273	1188	1191	1195	rVB	109730	156157	6.51%	0.350%
32	12.371	1195	1201	1205	rBV3	29416	76612	3.19%	0.172%
33	12.490	1205	1213	1215	rBV2	73643	166555	6.94%	0.373%
34	12.618	1222	1226	1229	rBV	1213250	1893856	78.97%	4.244%

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35	12.667	1229	1231	1235	rVV	365917	510800	21.30%	1.145%
36	12.786	1239	1243	1245	rBV2	276969	480966	20.06%	1.078%
37	12.825	1245	1247	1249	rVV	245400	370174	15.44%	0.830%
38	12.884	1249	1253	1258	rVB3	335164	741733	30.93%	1.662%
39	12.983	1260	1263	1265	rBV2	74298	120446	5.02%	0.270%
40	13.062	1268	1271	1275	rVB	195456	289890	12.09%	0.650%
41	13.150	1275	1280	1283	rBV2	220353	601813	25.09%	1.349%
42	13.200	1283	1285	1288	rBV	359612	622341	25.95%	1.395%
43	13.278	1291	1293	1298	rVV2	277095	576914	24.06%	1.293%
44	13.357	1298	1301	1306	rVB4	81663	202798	8.46%	0.455%
45	13.426	1306	1308	1310	rBV2	102142	140158	5.84%	0.314%
46	13.466	1310	1312	1317	rBV	135993	259010	10.80%	0.580%
47	13.545	1317	1320	1322	rBV	283883	356402	14.86%	0.799%
48	13.584	1322	1324	1326	rBV	463838	610627	25.46%	1.369%
49	13.623	1326	1328	1330	rVB2	152184	204399	8.52%	0.458%
50	13.663	1330	1332	1333	rBV	76562	105795	4.41%	0.237%
51	13.752	1338	1341	1344	rBV2	934884	1362193	56.80%	3.053%
52	13.801	1344	1346	1349	rVB3	187976	252775	10.54%	0.567%
53	13.850	1349	1351	1353	rBV	235274	314960	13.13%	0.706%
54	13.899	1353	1356	1360	rVV2	1283917	2255108	94.03%	5.054%
55	13.959	1360	1362	1364	rVV2	285370	408749	17.04%	0.916%
56	14.018	1364	1368	1371	rVV	811661	1206243	50.30%	2.703%
57	14.116	1376	1378	1381	rVV	460433	765320	31.91%	1.715%
58	14.195	1381	1386	1388	rVV3	564951	1358626	56.65%	3.045%
59	14.254	1388	1392	1398	rVV2	612247	1523534	63.53%	3.415%
60	14.343	1398	1401	1402	rVV2	104703	212113	8.84%	0.475%
61	14.402	1402	1407	1408	rVV3	138140	408608	17.04%	0.916%
62	14.451	1408	1412	1415	rVV2	547447	1332461	55.56%	2.986%
63	14.501	1415	1417	1419	rVV2	607612	984516	41.05%	2.206%
64	14.530	1419	1420	1424	rVV	301084	507099	21.14%	1.136%
65	14.599	1424	1427	1428	rVV	161781	299312	12.48%	0.671%
66	14.629	1428	1430	1433	rVV2	519671	794617	33.13%	1.781%
67	14.688	1433	1436	1437	rVV3	118885	251256	10.48%	0.563%
68	14.777	1441	1445	1448	rVV	1363121	2300717	95.93%	5.156%
69	14.856	1451	1453	1454	rVV	169931	247129	10.30%	0.554%
70	14.915	1455	1459	1463	rVV4	299530	961221	40.08%	2.154%
71	15.023	1467	1470	1474	rVV	670532	1131327	47.17%	2.536%

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72	15.092	1475	1477	1479	rVV3	96129	173099	7.22%	0.388%
73	15.141	1479	1482	1484	rVV	725294	1096767	45.73%	2.458%
74	15.191	1484	1487	1493	rVV4	211990	629226	26.24%	1.410%
75	15.270	1493	1495	1496	rVV	92183	126080	5.26%	0.283%
76	15.309	1496	1499	1502	rVV2	219124	461074	19.23%	1.033%
77	15.378	1502	1506	1509	rVV2	276901	688447	28.71%	1.543%
78	15.437	1509	1512	1519	rVV	167554	398557	16.62%	0.893%
79	15.536	1519	1522	1523	rVV3	48270	105443	4.40%	0.236%
80	15.555	1523	1524	1527	rVV2	63909	82572	3.44%	0.185%
81	15.624	1527	1531	1532	rVV4	23253	46523	1.94%	0.104%
82	15.674	1532	1536	1541	rVV	253764	495067	20.64%	1.110%
83	15.743	1541	1543	1549	rVB4	50618	115282	4.81%	0.258%
84	15.861	1549	1555	1558	rBV7	55003	112841	4.71%	0.253%
85	15.969	1563	1566	1568	rBV4	47465	97427	4.06%	0.218%
86	16.029	1568	1572	1578	rVB3	48216	168447	7.02%	0.378%

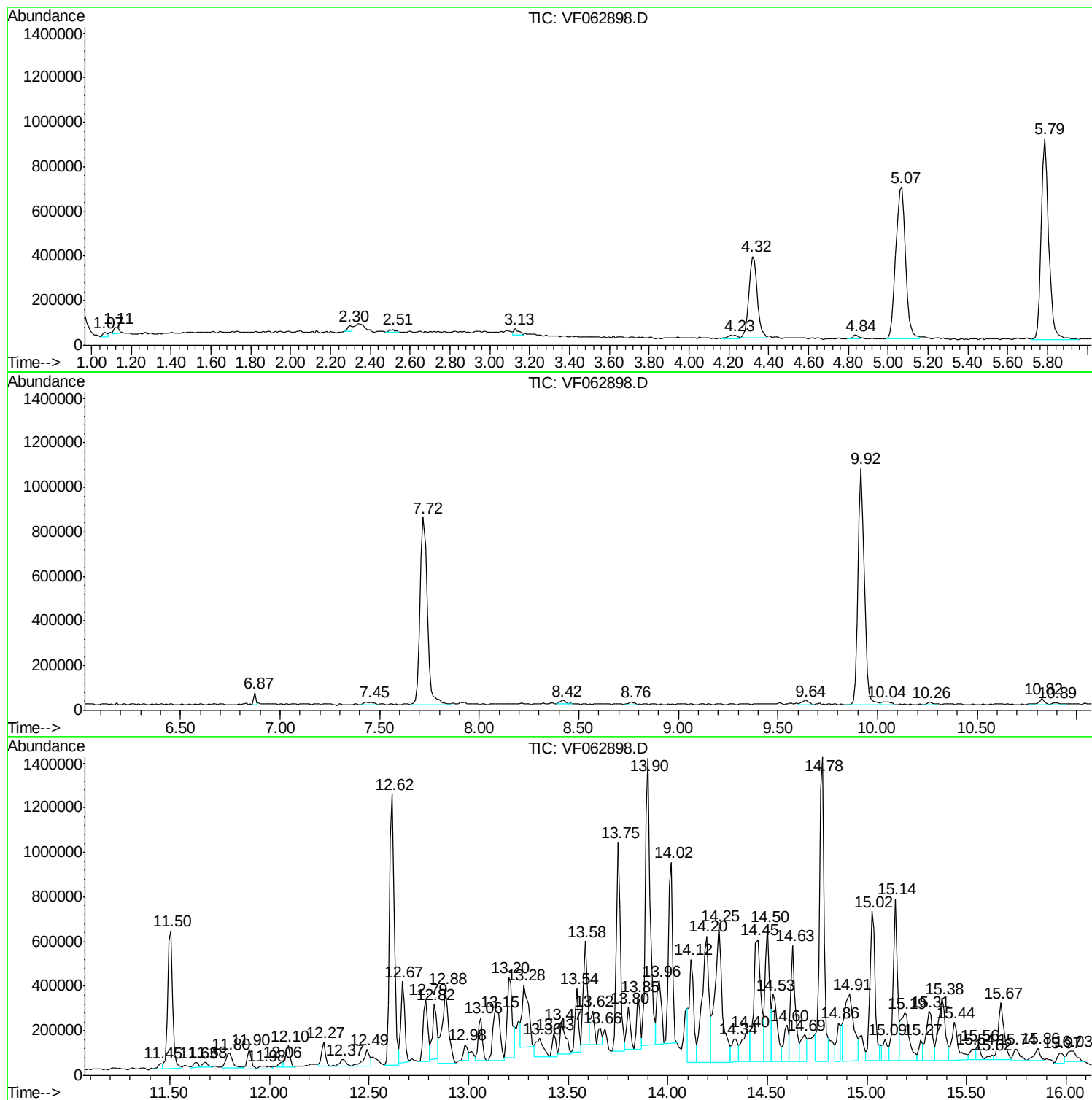
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P



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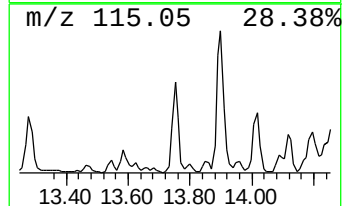
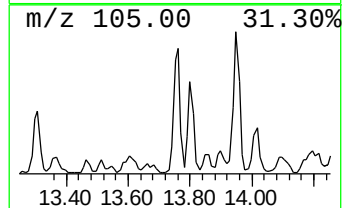
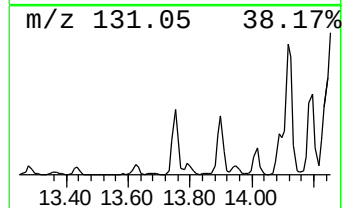
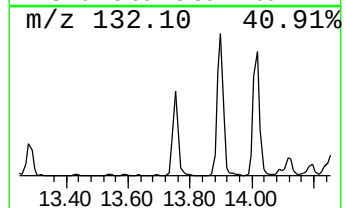
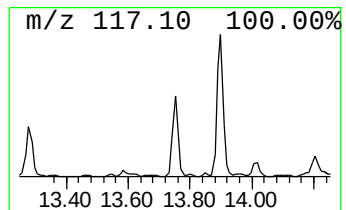
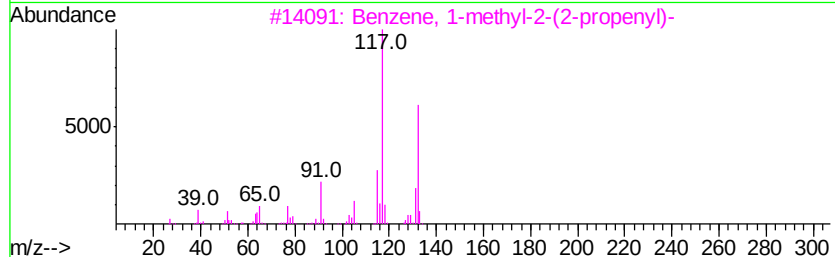
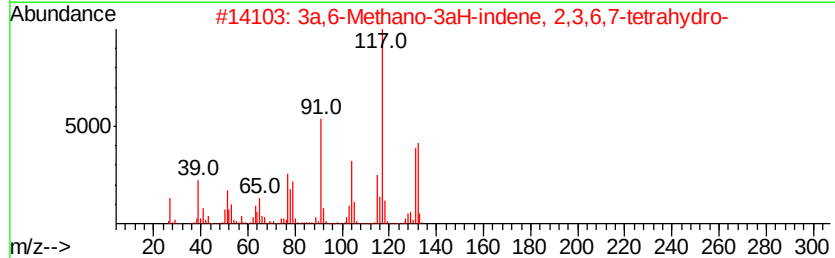
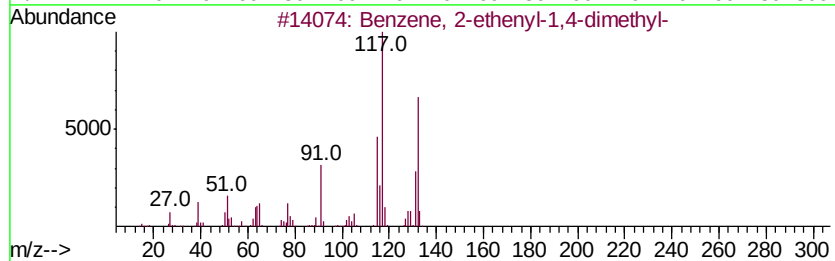
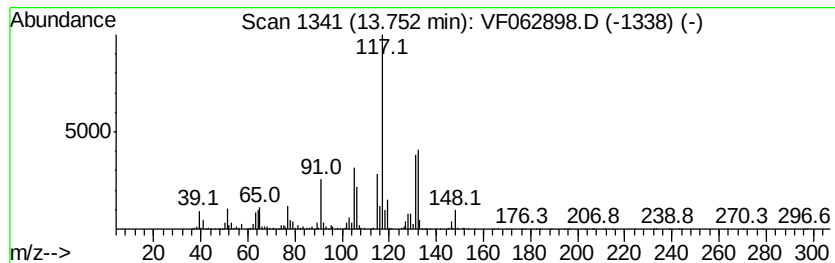
Instrument :
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ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.75	35.96 ug/l	1362190	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 2-ethenyl-1,4-dimethyl-	132 C10H12	002039-89-6 89
2		3a,6-Methano-3aH-indene, 2,3,6,7...	132 C10H12	098640-29-0 76
3		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 70
4		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 64
5		Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7 64



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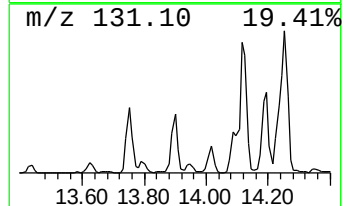
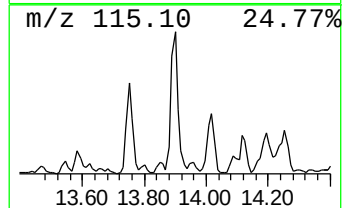
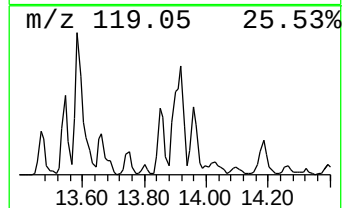
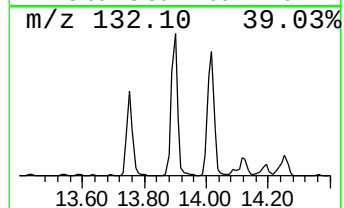
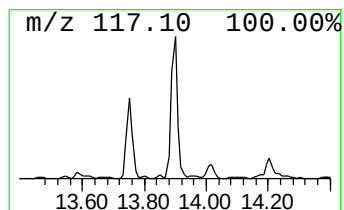
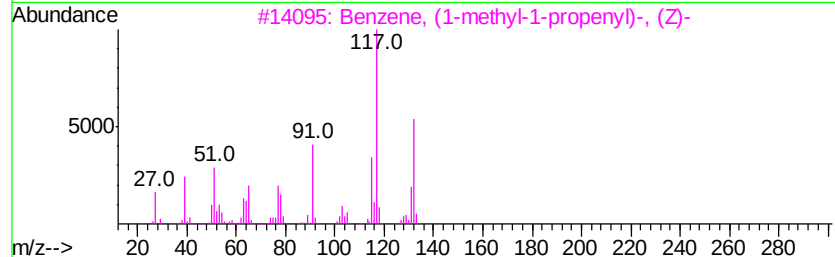
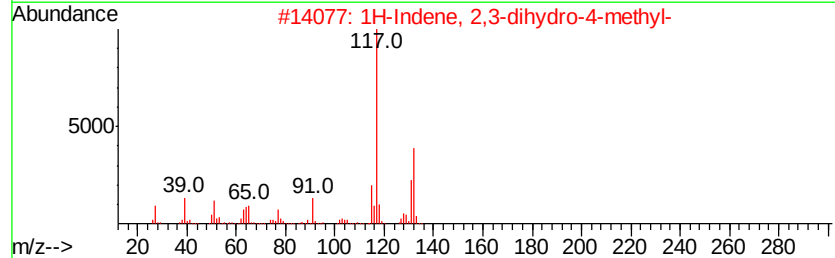
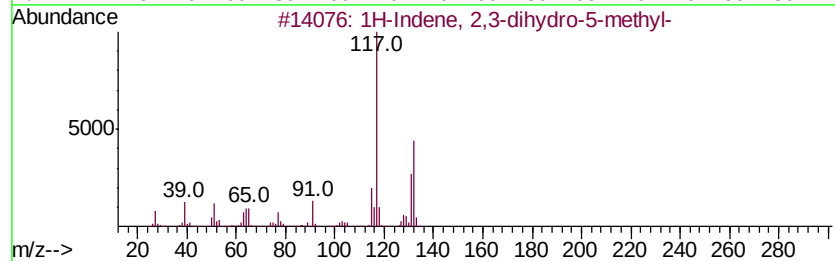
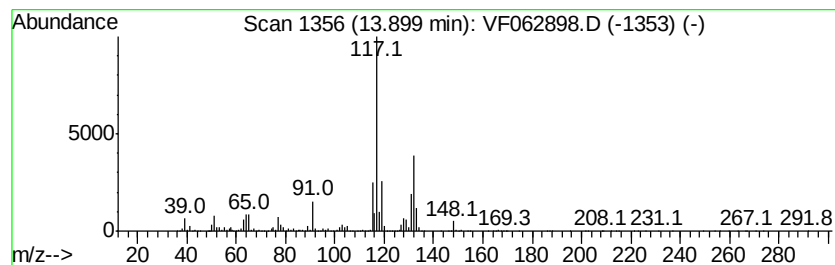
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	59.54 ug/l	2255110	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-5-methyl-	132 C10H12	000874-35-1	94
2	1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6	91
3	Benzene, (1-methyl-1-propenyl)-,...	132 C10H12	000767-99-7	80
4	Benzene, 1-ethenyl-4-ethyl-	132 C10H12	003454-07-7	76
5	1-Phenyl-1-butene	132 C10H12	000824-90-8	74



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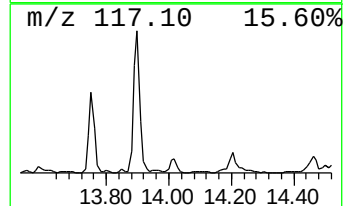
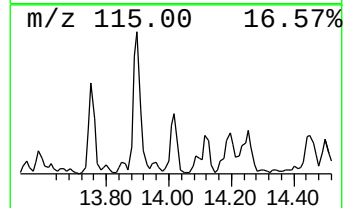
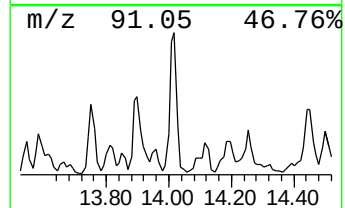
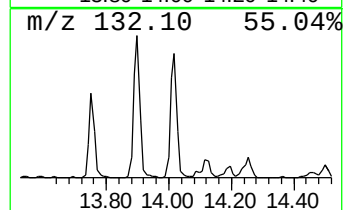
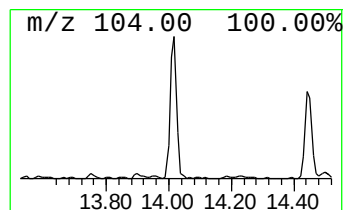
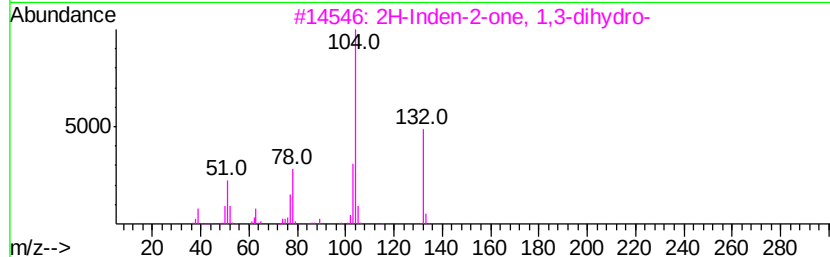
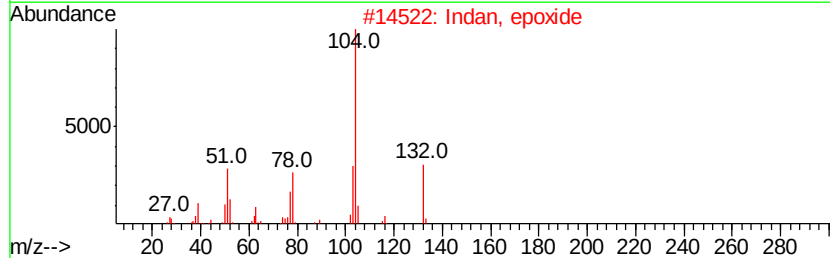
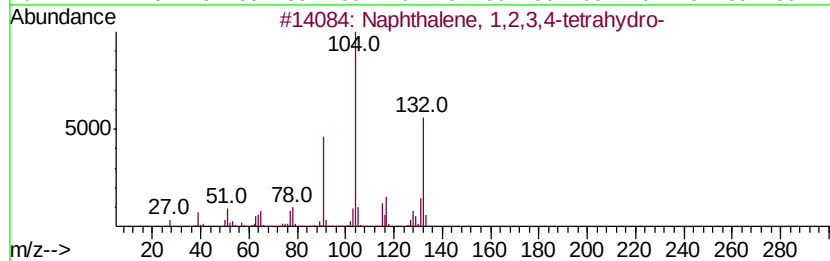
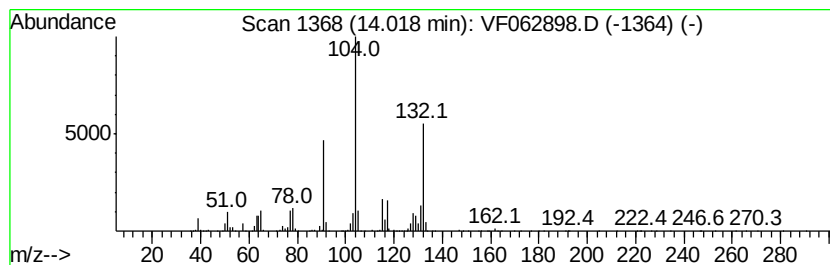
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 3 Naphthalene, 1,2,3,4-tetrahydro- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.02	31.85 ug/l	1206240	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-	132 C10H12	000119-64-2 96
2		Indan, epoxide	132 C9H8O	000768-22-9 58
3		2H-Inden-2-one, 1,3-dihydro-	132 C9H8O	000615-13-4 53
4		Naphthalene, 1,4,5,8-tetrahydro-	132 C10H12	000493-04-9 41
5		2-Propenoic acid, 2-phenylethyl ...	176 C11H12O2	003530-36-7 38



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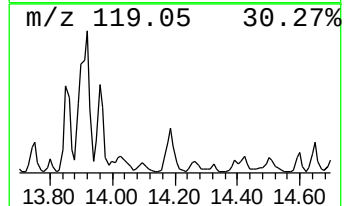
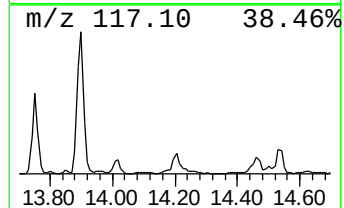
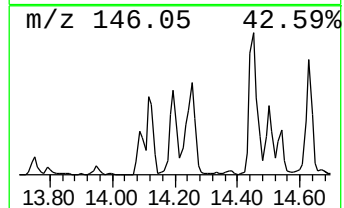
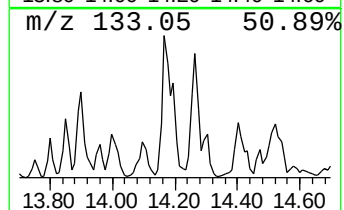
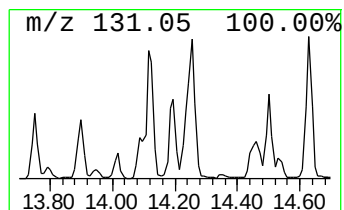
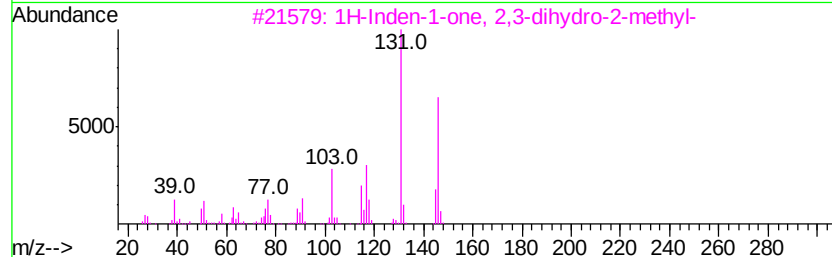
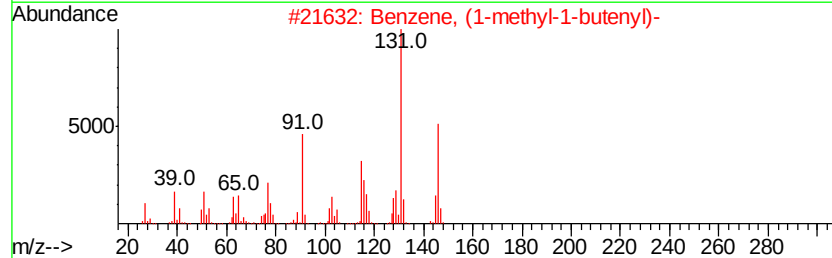
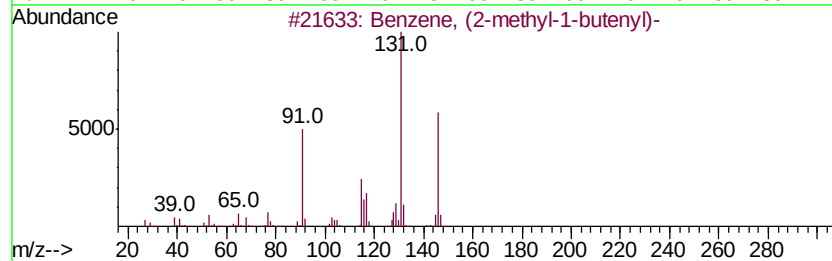
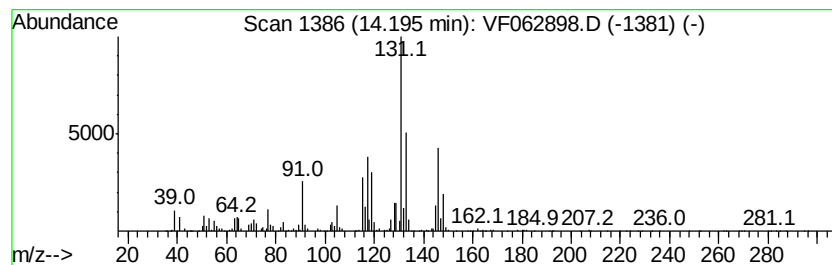
Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 4 Benzene, (2-methyl-1-butenyl)- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.20	35.87 ug/l	1358630	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 90
2		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 78
3		1H-Inden-1-one, 2,3-dihydro-2-me...	146 C10H10O	017496-14-9 64
4		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 55
5		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 55



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

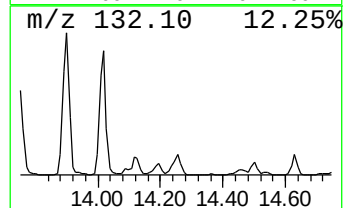
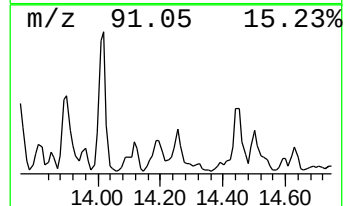
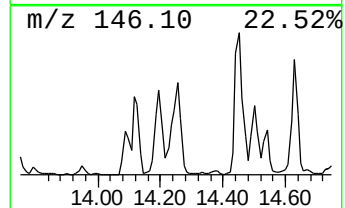
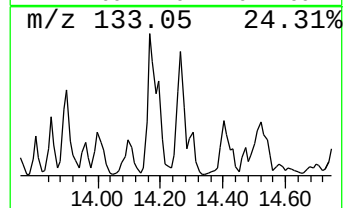
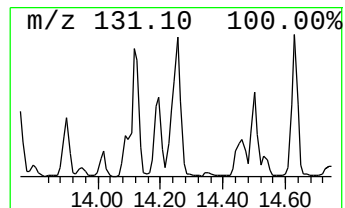
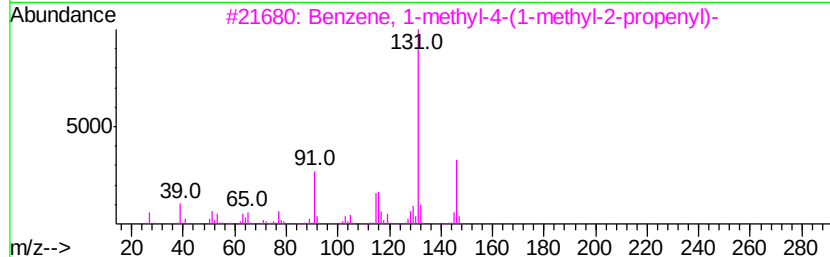
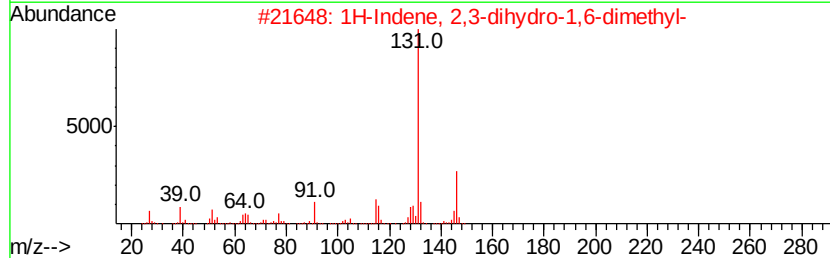
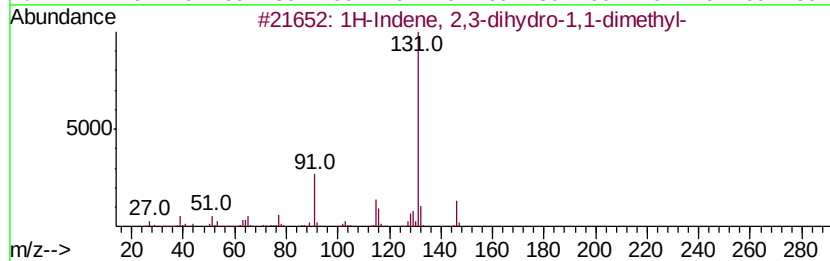
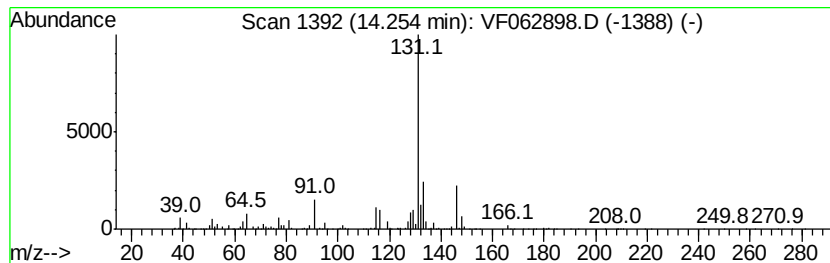
Instrument :
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ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.25	40.22 ug/l	1523530	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9	81
2	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	81
3	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	81
4	Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5	74
5	2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

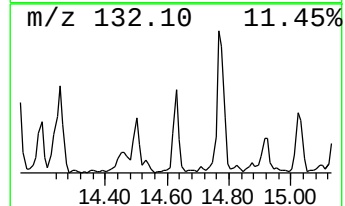
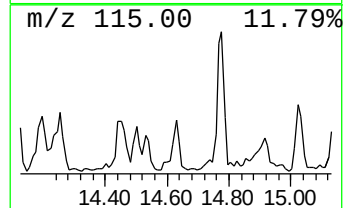
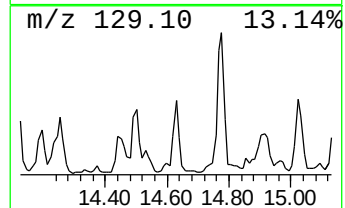
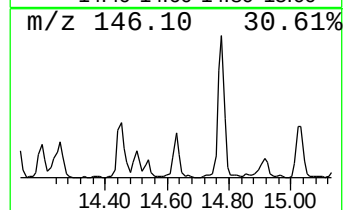
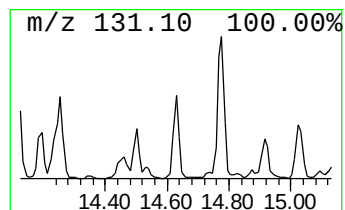
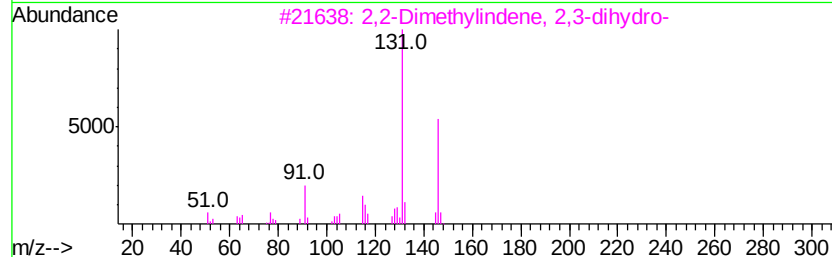
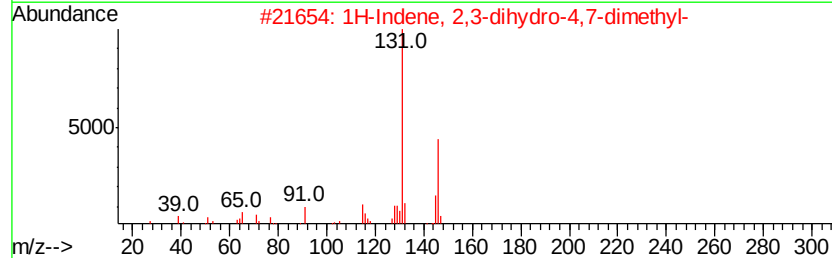
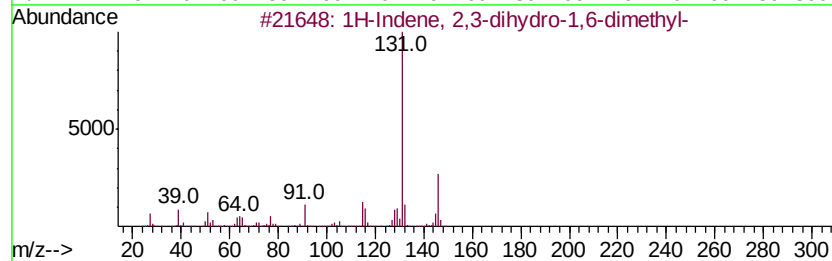
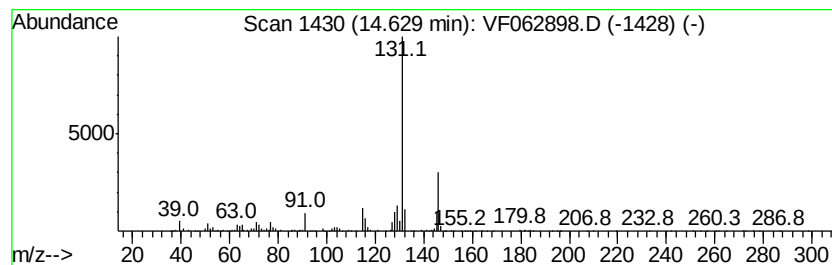
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Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 6 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.63	20.98 ug/l	794617	1,4-Dichlorobenzene-d4	12.62

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	96
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	93
3		2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	91
4		1H-Indene, 2,3-dihydro-1,2-dimet...	146	C11H14	017057-82-8	91
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	91



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

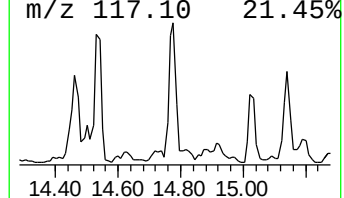
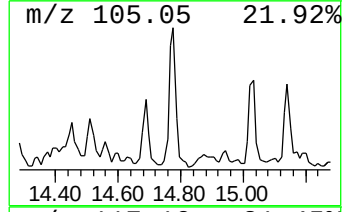
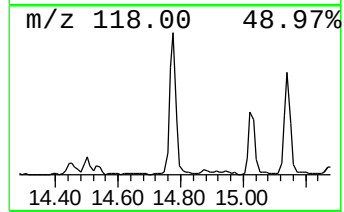
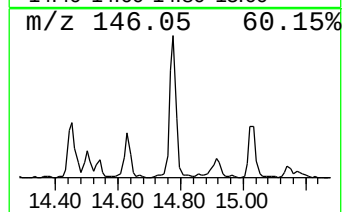
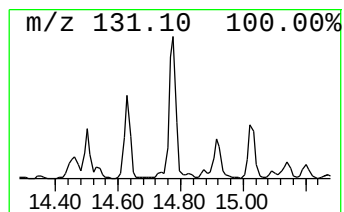
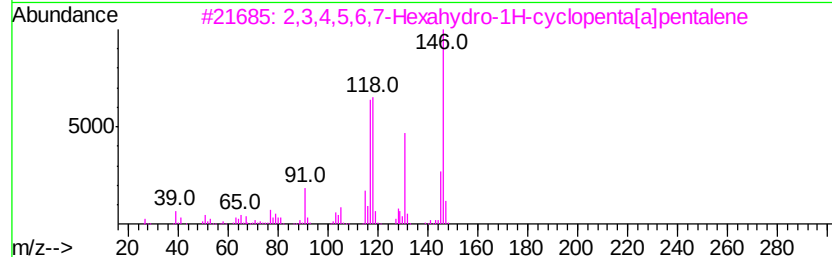
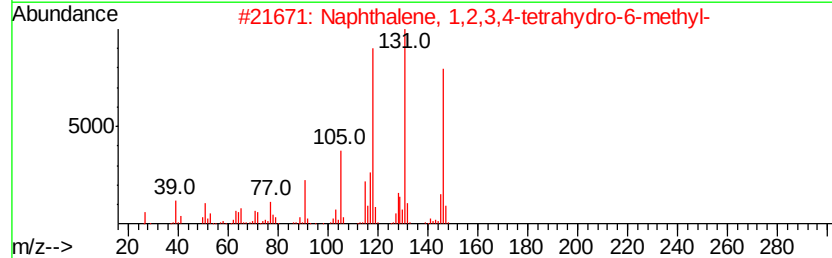
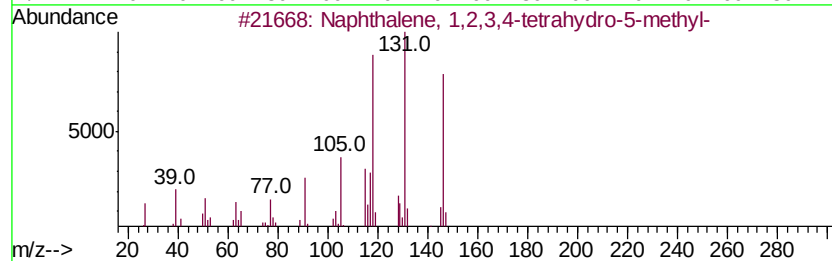
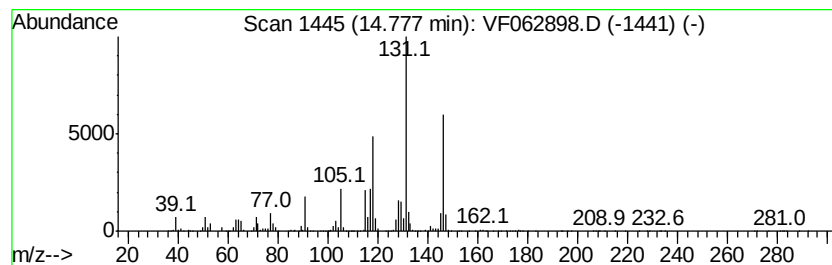
Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 7 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.78	60.74 ug/l	2300720	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 96
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 96
3		2,3,4,5,6,7-Hexahydro-1H-cyclope...	146 C11H14	1000189-31-0 76
4		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 70
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 70



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

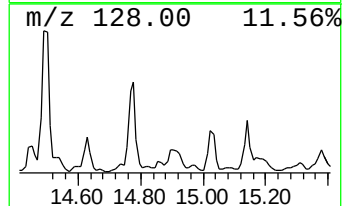
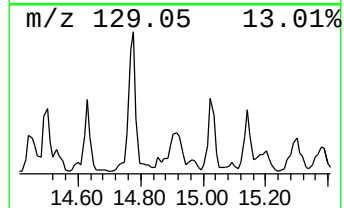
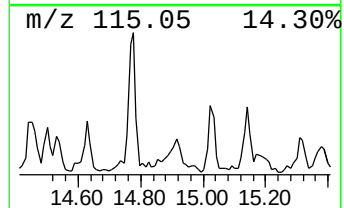
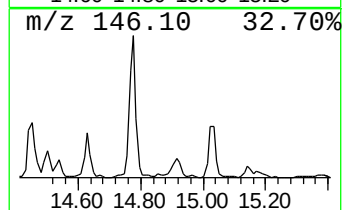
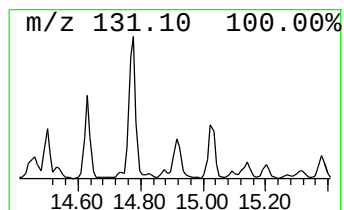
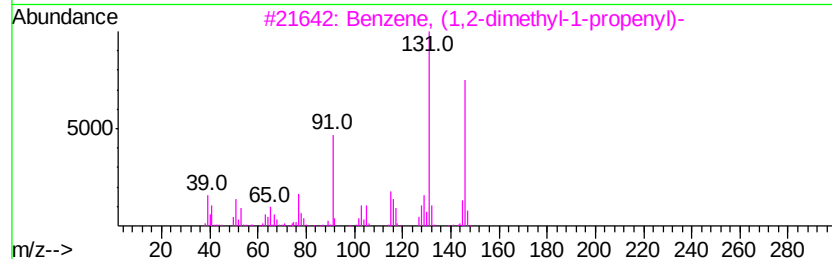
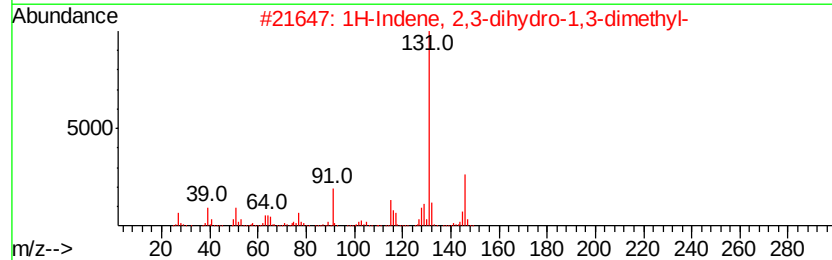
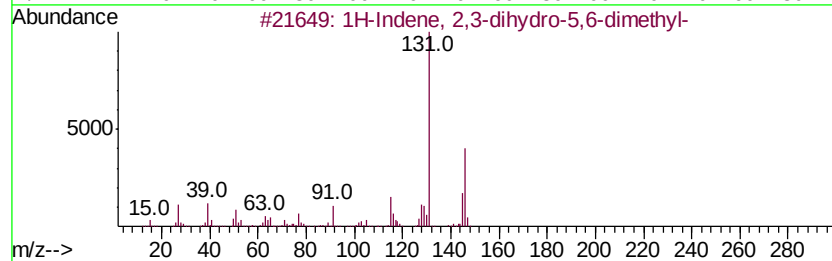
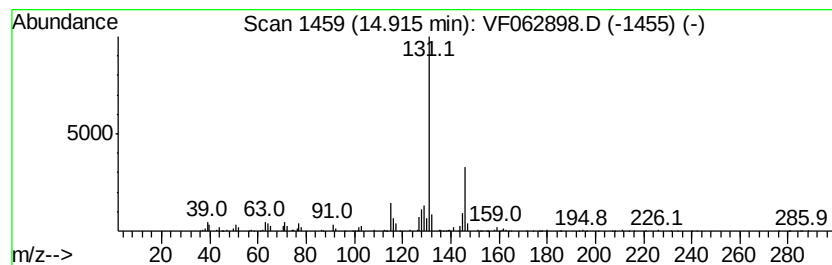
Instrument :
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ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 8 1H-Indene, 2,3-dihydro-5,6-... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.91	25.38 ug/l	961221	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	1H-Indene, 2,3-dihydro-5,6-dimet...	146 C11H14	001075-22-5	93
2	1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5	91
3	Benzene, (1,2-dimethyl-1-propenyl)-	146 C11H14	000769-57-3	91
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9	91
5	1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

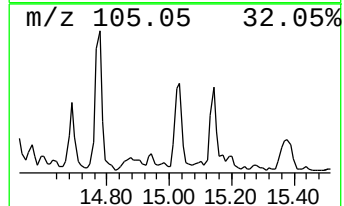
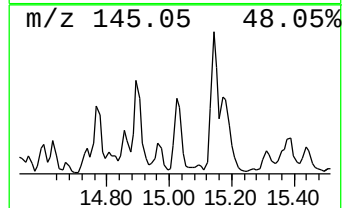
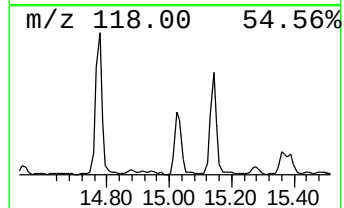
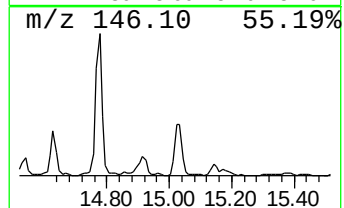
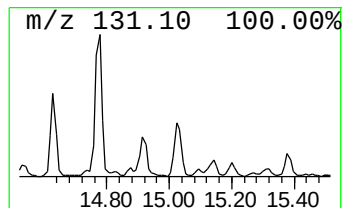
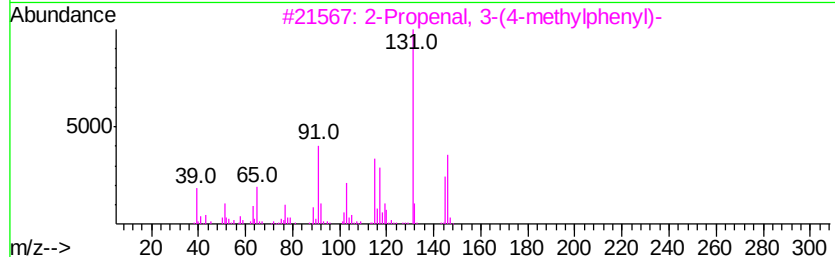
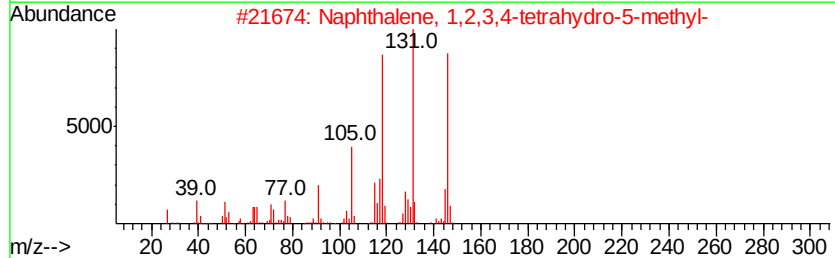
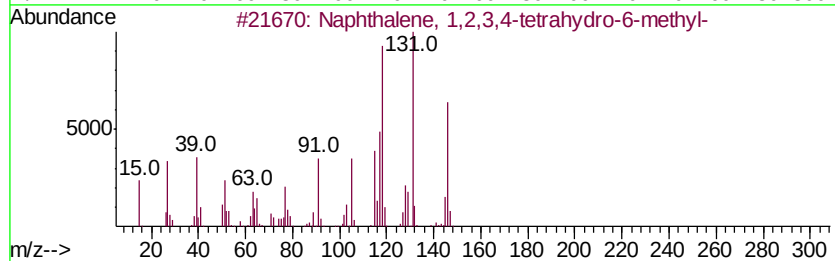
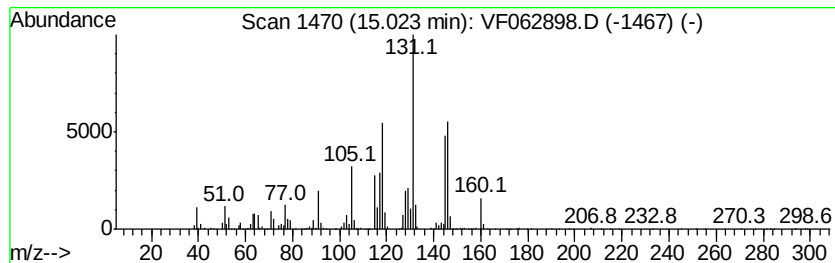
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ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 9 Naphthalene, 1,2,3,4-tetrahydro-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	29.87 ug/l	1131330	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 93
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 81
3		2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2 64
4		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 64
5		1H-1,3,2-Benzodiazaborole, 2-eth...	146 C8H11BN2	074663-81-3 62



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
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Sample : K3160-01
Misc : 7.75g/5mL/MSVOA F/SOIL
ALS Vial : 7 Sample Multiplier: 1

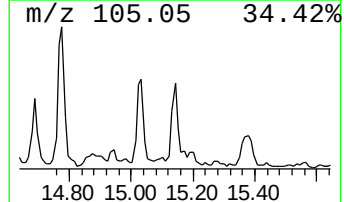
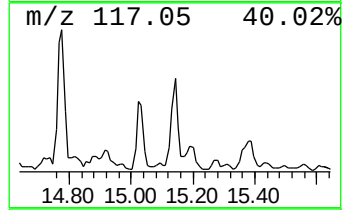
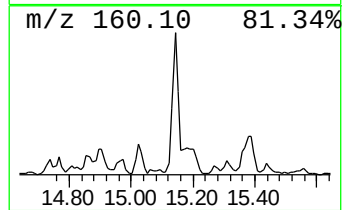
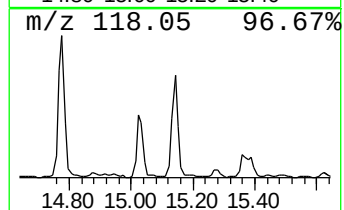
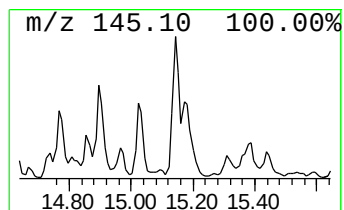
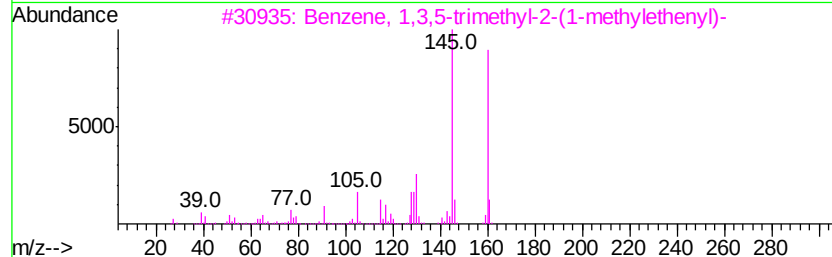
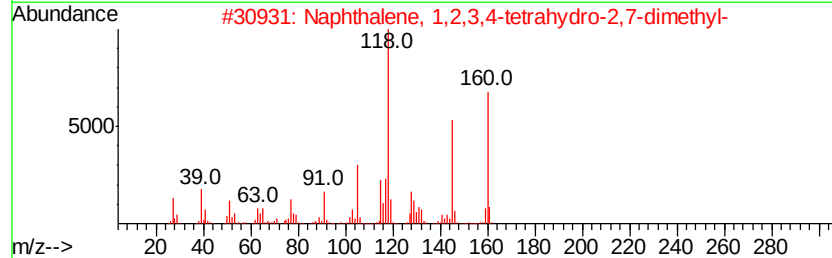
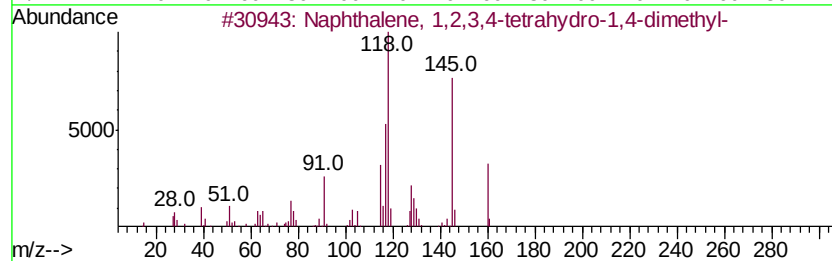
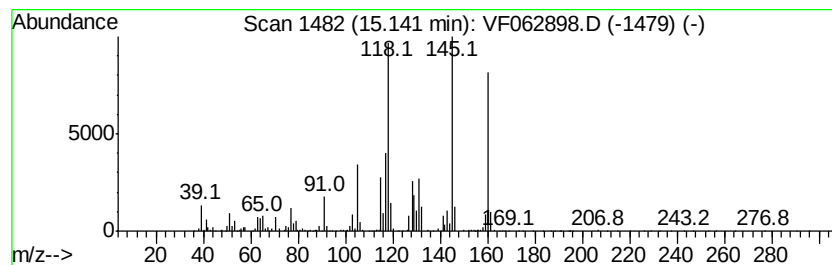
Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.14	28.96 ug/l	1096770	1,4-Dichlorobenzene-d4	12.62
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	004175-54-6 95
2		Naphthalene, 1,2,3,4-tetrahydro-...	160 C12H16	013065-07-1 95
3		Benzene, 1,3,5-trimethyl-2-(1-me...	160 C12H16	014679-13-1 78
4		Benzene, 4-(2-butenyl)-1,2-dimet...	160 C12H16	054340-86-2 64
5		Benzene, 1-(1-methylethenyl)-3-(...	160 C12H16	001129-29-9 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA_F\DATA\VF060719\
Data File : VF062898.D
Acq On : 7 Jun 2019 14:23
Operator : FY/SY
Sample : K3160-01
Misc : 7.75g/5mL/MSVOA_F/SOIL
ALS Vial : 7 Sample Multiplier: 1

Instrument :
MSVOA_F
ClientSampleId :
EO-02-060619-A

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_F\METHODS\82F053019S.M
Quant Title : SW846 8260

TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 2-etheny...	13.75	36.0	ug/l	1362190	4	12.62	1893860	50.0
1H-Indene, 2,3-di...	13.90	59.5	ug/l	2255110	4	12.62	1893860	50.0
Naphthalene, 1,2,...	14.02	31.9	ug/l	1206240	4	12.62	1893860	50.0
Benzene, (2-methy...	14.20	35.9	ug/l	1358630	4	12.62	1893860	50.0
1H-Indene, 2,3-di...	14.25	40.2	ug/l	1523530	4	12.62	1893860	50.0
1H-Indene, 2,3-di...	14.63	21.0	ug/l	794617	4	12.62	1893860	50.0
Naphthalene, 1,2,...	14.78	60.7	ug/l	2300720	4	12.62	1893860	50.0
1H-Indene, 2,3-di...	14.91	25.4	ug/l	961221	4	12.62	1893860	50.0
Naphthalene, 1,2,...	15.02	29.9	ug/l	1131330	4	12.62	1893860	50.0
Naphthalene, 1,2,...	15.14	29.0	ug/l	1096770	4	12.62	1893860	50.0