

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD120420\
 Data File : VD067880.D
 Acq On : 04 Dec 2020 12:30
 Operator : VA/SY
 Sample : L4939-01
 Misc : 7.46G/5.00ml/MSVOA D/SOIL
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 EO-02-12320

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_D\METHOD\82D120120S.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	4.973	509	519	526	rBV5	6202	17958	1.68%	0.166%
2	7.920	1010	1020	1025	rBV2	145834	350518	32.81%	3.249%
3	7.985	1025	1031	1042	rVB2	297476	677079	63.39%	6.276%
4	8.338	1080	1091	1102	rVB	130010	313206	29.32%	2.903%
5	8.867	1171	1181	1191	rBV	351346	723940	67.77%	6.711%
6	9.356	1257	1264	1272	rVB4	11189	24176	2.26%	0.224%
7	10.344	1423	1432	1448	rBV	568705	1068196	100.00%	9.902%
8	10.514	1454	1461	1469	rVB6	15077	34928	3.27%	0.324%
9	10.685	1482	1490	1496	rBV3	33721	61303	5.74%	0.568%
10	10.779	1501	1506	1518	rVB5	15802	35352	3.31%	0.328%
11	11.197	1572	1577	1582	rVB3	23369	38145	3.57%	0.354%
12	11.250	1582	1586	1592	rVB4	13657	23284	2.18%	0.216%
13	11.461	1618	1622	1630	rVB3	9302	15330	1.44%	0.142%
14	11.650	1645	1654	1665	rVB	489657	839445	78.59%	7.782%
15	11.744	1665	1670	1674	rBV4	9247	17083	1.60%	0.158%
16	11.891	1693	1695	1700	rVB4	12641	16408	1.54%	0.152%
17	12.155	1733	1740	1742	rBV5	16244	27837	2.61%	0.258%
18	12.273	1754	1760	1764	rVB6	8575	15918	1.49%	0.148%
19	12.355	1770	1774	1777	rBV5	9996	17056	1.60%	0.158%
20	12.638	1816	1822	1833	rVB	403720	683356	63.97%	6.335%
21	12.802	1845	1850	1857	rVB6	21773	44721	4.19%	0.415%
22	12.879	1857	1863	1864	rBV4	8815	15283	1.43%	0.142%
23	12.961	1866	1877	1882	rVV4	97925	237107	22.20%	2.198%
24	13.014	1882	1886	1891	rVB7	18062	32547	3.05%	0.302%
25	13.126	1897	1905	1910	rBV4	33427	81715	7.65%	0.757%
26	13.191	1912	1916	1928	rVV6	32142	74593	6.98%	0.691%
27	13.402	1948	1952	1955	rVB4	11278	16197	1.52%	0.150%
28	13.467	1956	1963	1967	rVV6	52565	105696	9.89%	0.980%
29	13.514	1967	1971	1975	rVV5	34238	68158	6.38%	0.632%
30	13.579	1976	1982	1991	rVV2	510294	1005899	94.17%	9.325%
31	13.649	1991	1994	1998	rVV4	14158	18477	1.73%	0.171%
32	13.720	2001	2006	2007	rVV5	11315	15405	1.44%	0.143%
33	13.738	2007	2009	2012	rVV2	19843	29031	2.72%	0.269%
34	13.773	2012	2015	2018	rVV3	30442	50369	4.72%	0.467%

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Integration Parameters: RTEINT.P

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

Filtering: 5
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Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
 Title : SW846 8260

35	13.814	2018	2022	2031	rVB2	112465	185456	17.36%	1.719%
36	13.902	2031	2037	2042	rBV2	153889	254361	23.81%	2.358%
37	13.973	2045	2049	2059	rVB	52182	98438	9.22%	0.913%
38	14.061	2059	2064	2066	rBV6	11574	18860	1.77%	0.175%
39	14.120	2069	2074	2079	rVB	80677	119812	11.22%	1.111%
40	14.232	2087	2093	2097	rVB3	96429	157897	14.78%	1.464%
41	14.279	2097	2101	2109	rVB8	23394	58354	5.46%	0.541%
42	14.367	2109	2116	2120	rBV2	60925	116616	10.92%	1.081%
43	14.414	2120	2124	2126	rVV3	61416	95511	8.94%	0.885%
44	14.443	2127	2129	2133	rVV2	52442	80669	7.55%	0.748%
45	14.485	2133	2136	2139	rVV	64745	77076	7.22%	0.714%
46	14.526	2139	2143	2147	rVV2	59804	81220	7.60%	0.753%
47	14.573	2147	2151	2158	rVB4	64558	117978	11.04%	1.094%
48	14.667	2158	2167	2172	rVB5	47596	98794	9.25%	0.916%
49	14.714	2172	2175	2177	rBV3	21714	28254	2.65%	0.262%
50	14.755	2178	2182	2187	rVB3	72515	117165	10.97%	1.086%
51	14.820	2187	2193	2200	rBV3	182128	418680	39.20%	3.881%
52	14.885	2200	2204	2212	rVB3	48634	87303	8.17%	0.809%
53	14.991	2217	2222	2227	rVB2	56911	92194	8.63%	0.855%
54	15.102	2236	2241	2244	rBV3	66976	111539	10.44%	1.034%
55	15.214	2254	2260	2266	rVB3	92310	188603	17.66%	1.748%
56	15.367	2280	2286	2290	rBV6	32450	66579	6.23%	0.617%
57	15.461	2296	2302	2307	rBV2	70462	121631	11.39%	1.128%
58	15.514	2307	2311	2312	rVV3	43299	59111	5.53%	0.548%
59	15.538	2313	2315	2323	rVV4	52476	123452	11.56%	1.144%
60	15.602	2323	2326	2335	rVB7	41311	79340	7.43%	0.735%
61	15.702	2335	2343	2351	rBV6	33706	102877	9.63%	0.954%
62	15.779	2351	2356	2360	rBV7	16440	32242	3.02%	0.299%
63	15.867	2362	2371	2375	rBV7	36155	101276	9.48%	0.939%
64	15.908	2375	2378	2384	rVB3	57144	95549	8.94%	0.886%
65	16.002	2389	2394	2399	rVV8	21168	42409	3.97%	0.393%
66	16.079	2400	2407	2422	rVB5	36323	147743	13.83%	1.370%
67	16.232	2423	2433	2444	rBV4	68718	184174	17.24%	1.707%
68	16.432	2458	2467	2474	rBV5	65438	162721	15.23%	1.508%
69	16.502	2474	2479	2484	rVB4	18503	35175	3.29%	0.326%
70	16.696	2505	2512	2513	rBV4	20536	32823	3.07%	0.304%

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Integration Parameters: RTEINT.P
Integrator: RTE
Smoothing : ON Filtering: 5
Sampling : 1 Min Area: 3 % 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:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Title : SW846 8260

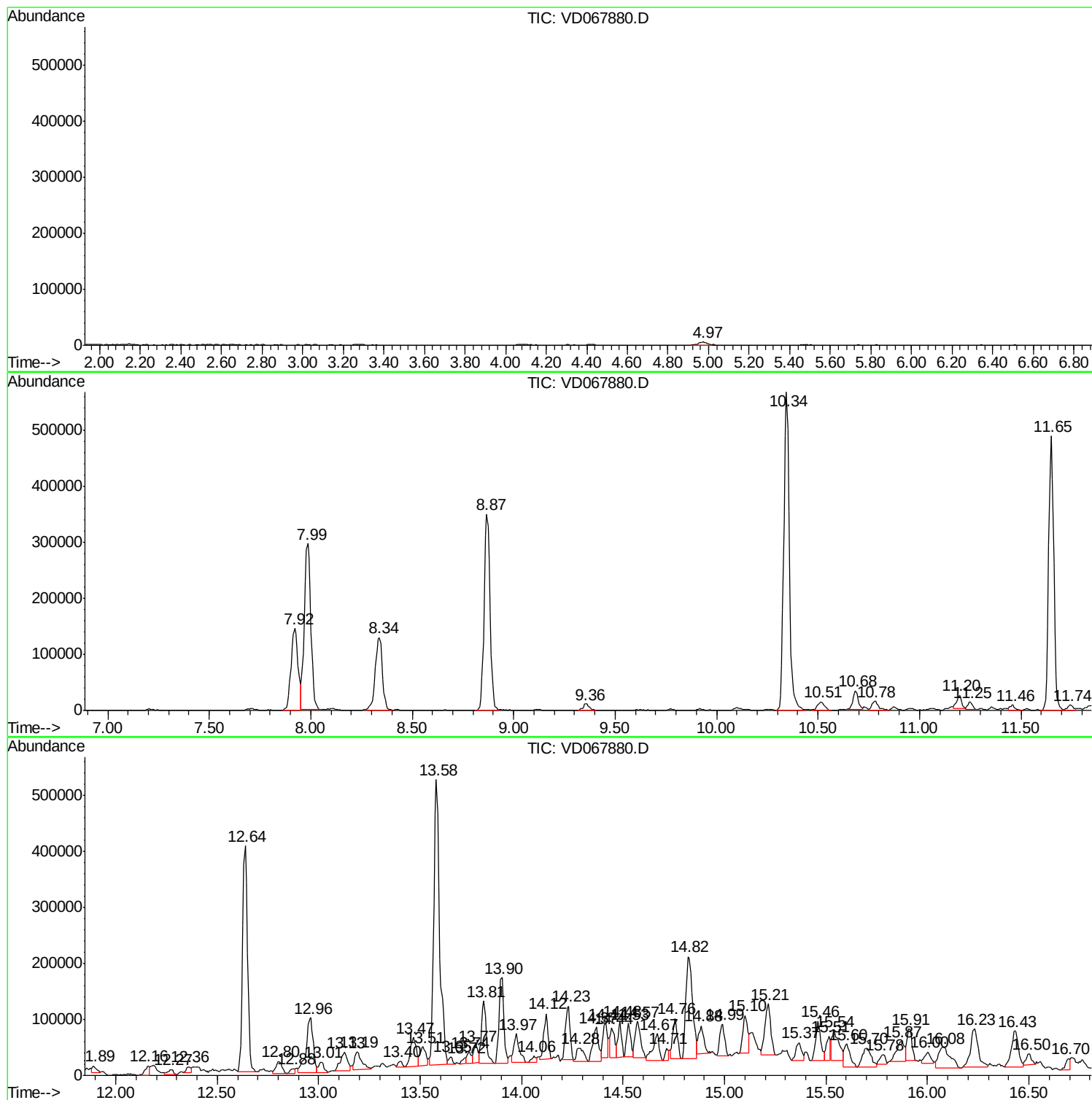
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Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
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Instrument :
MSVOA_D
ClientSampleId :
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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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



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Operator : VA/SY
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Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

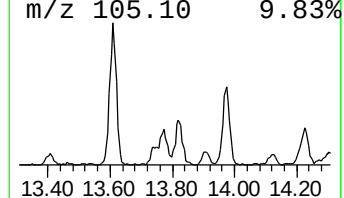
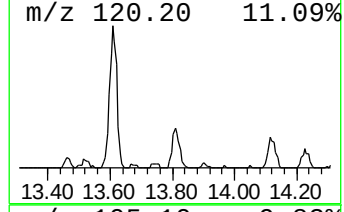
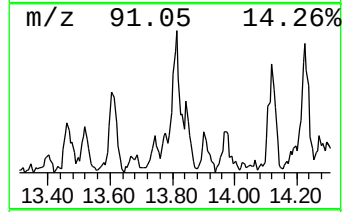
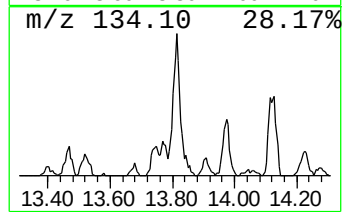
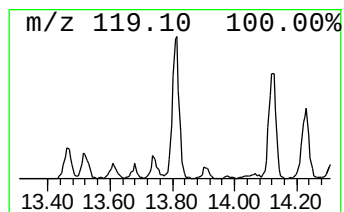
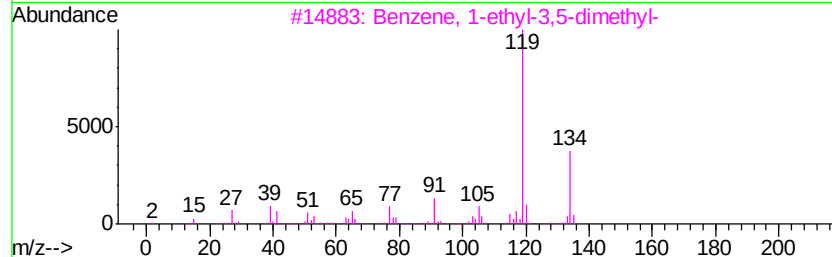
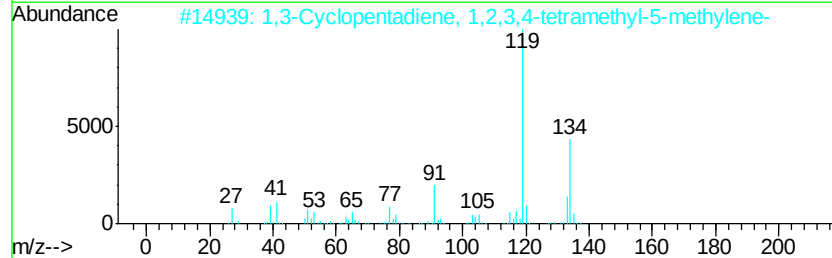
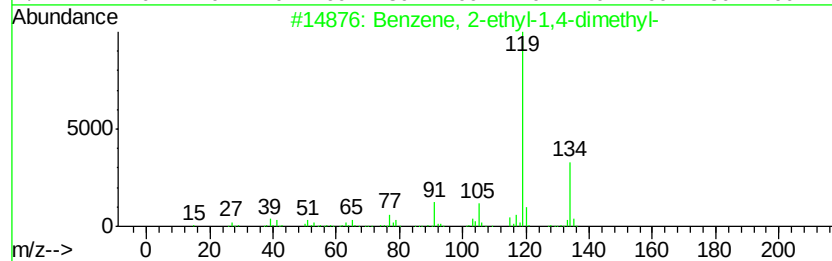
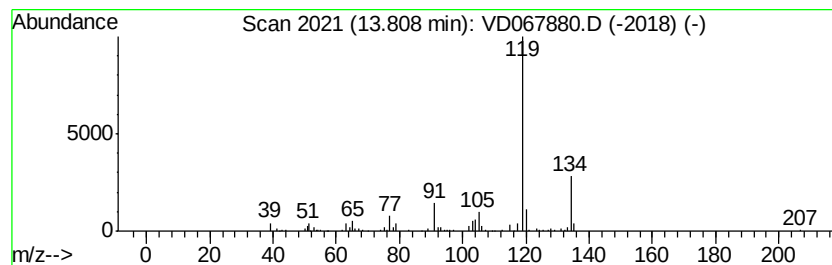
Instrument :
MSVOA_D
ClientSampled :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.		
13.81	9.22 ug/l	185456	1,4-Dichlorobenzene-d4	13.58		
Hit# of	5	Tentative ID	MW	MolForm	CAS#	Qual
1		Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87
2		1,3-Cyclopentadiene, 1,2,3,4-tet...	134	C10H14	076089-59-3	87
3		Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	87
4		Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	87
5		Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	86



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Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

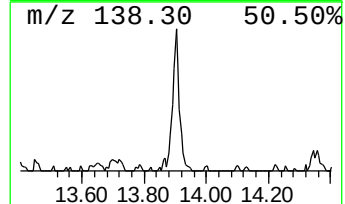
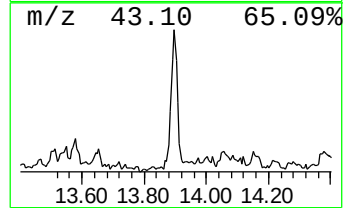
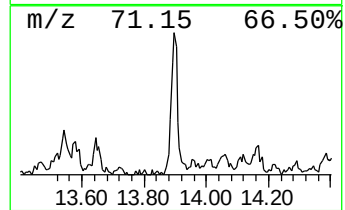
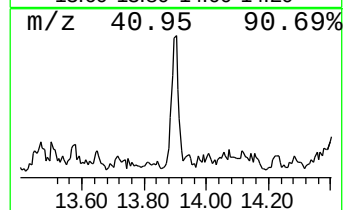
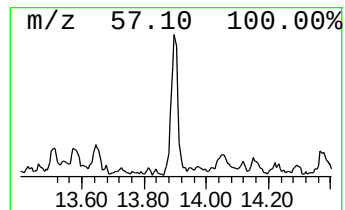
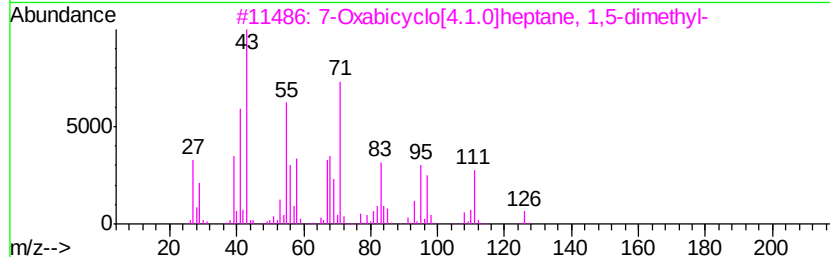
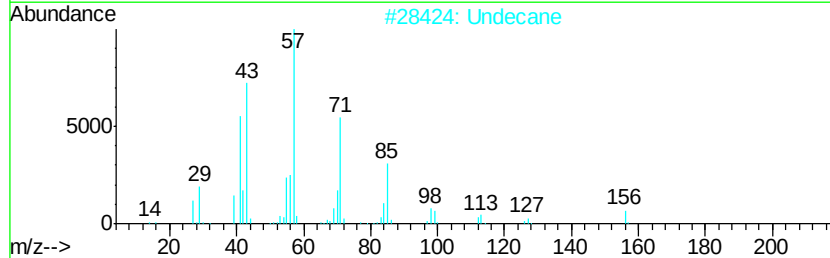
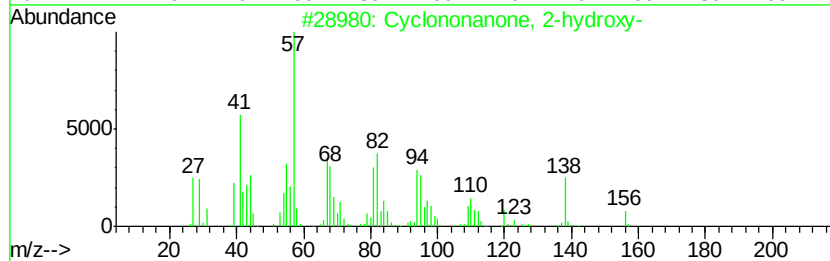
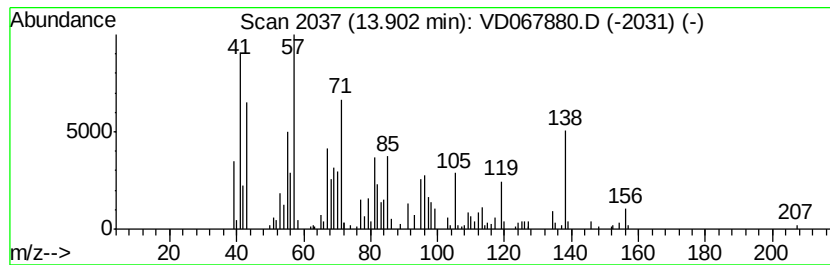
Instrument :
MSVOA_D
ClientSampleId :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 2 unknown13.90 Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.	
13.90	12.64 ug/l	254361	1,4-Dichlorobenzene-d4	13.58	
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclononane, 2-hydroxy-	156	C9H16O2	000496-83-3	38
2	Undecane	156	C11H24	001120-21-4	35
3	7-Oxabicyclo[4.1.0]heptane, 1,5-...	126	C8H14O	162239-52-3	30
4	Citronellol	156	C10H20O	000106-22-9	25
5	4-Acetyl-1-methylcyclohexene	138	C9H14O	006090-09-1	25



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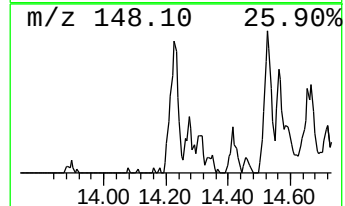
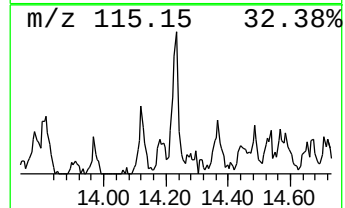
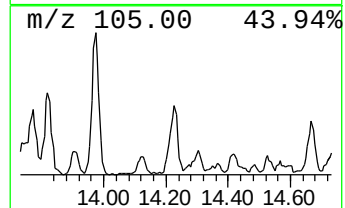
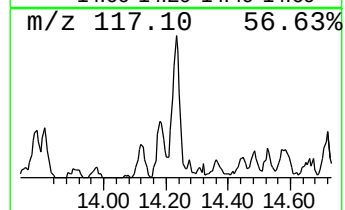
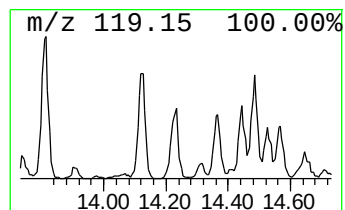
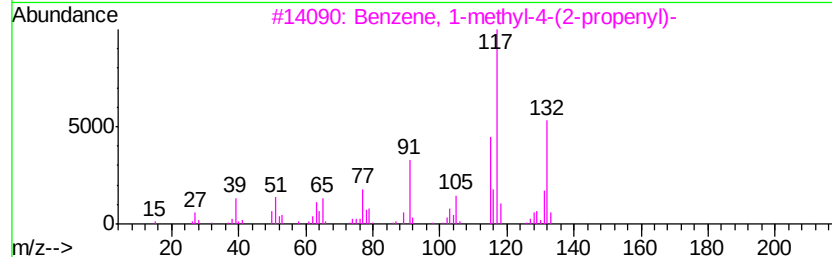
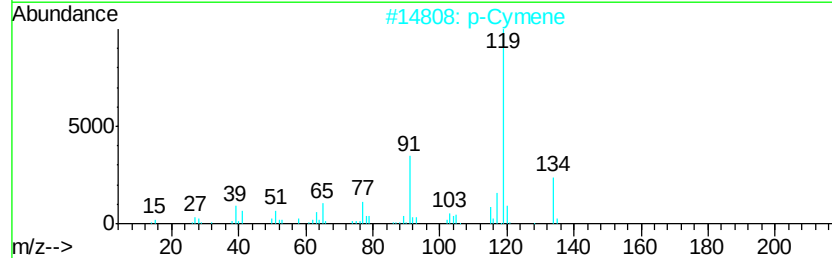
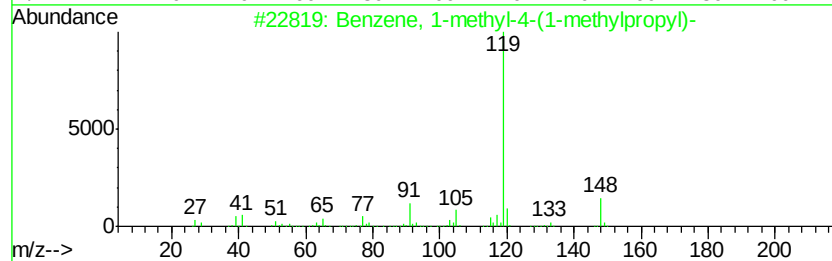
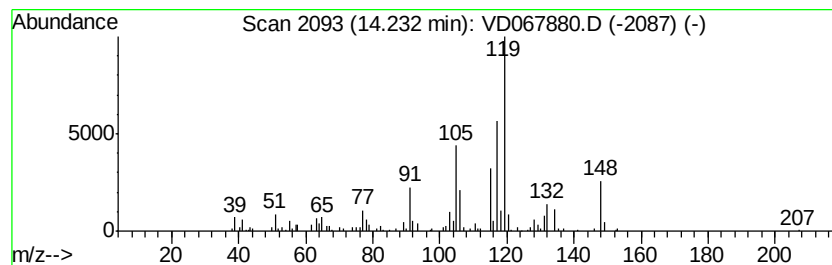
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 3 unknown14.23 Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.23	7.85 ug/l	157897	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	46
2		p-Cymene	134	C10H14	000099-87-6	45
3		Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	42
4		4'-Methylpropiophenone	148	C10H12O	005337-93-9	38
5		Benzene, 1,3-diethyl-5-methyl-	148	C11H16	002050-24-0	38



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ALS Vial : 6 Sample Multiplier: 1

Instrument :
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ClientSampleId :
EO-02-12320

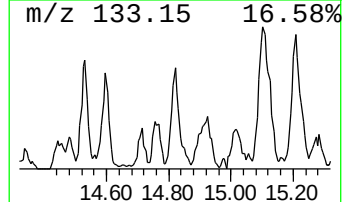
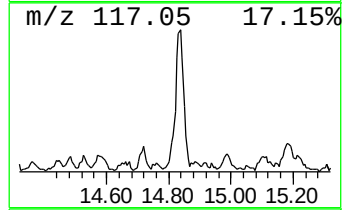
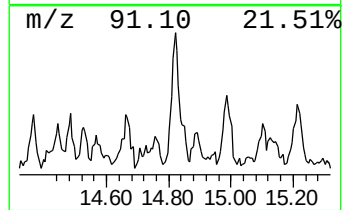
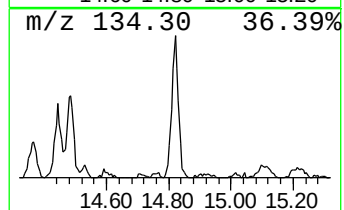
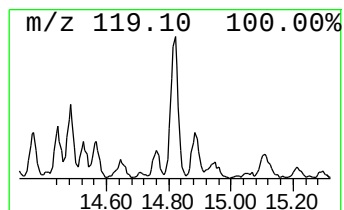
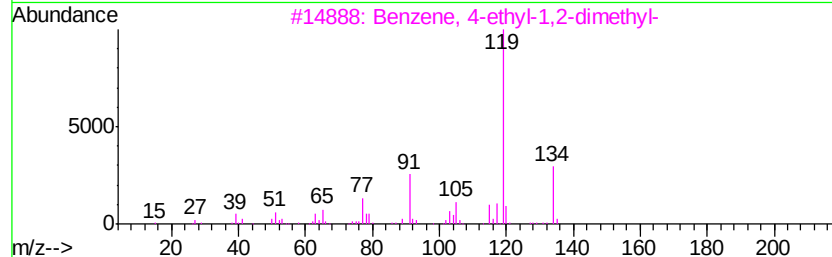
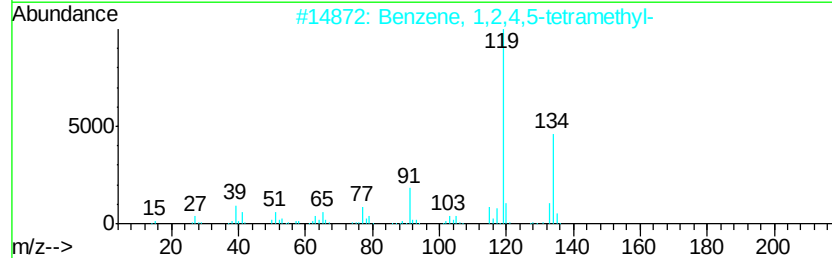
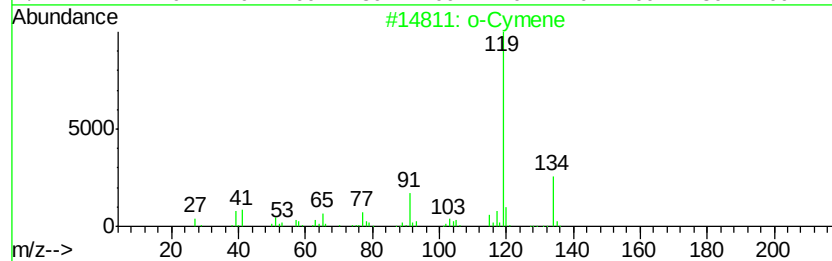
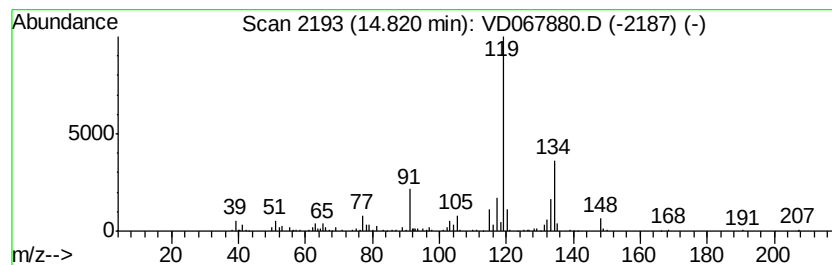
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Quant Title : SW846 8260

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

Peak Number 4 o-Cymene Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.82	20.81 ug/l	418680	1,4-Dichlorobenzene-d4	13.58

Hit# of	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	94
2	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	93
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	91
4	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	87
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	87



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

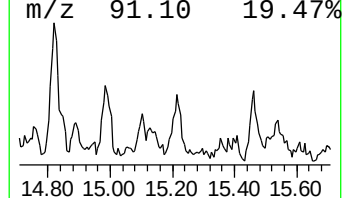
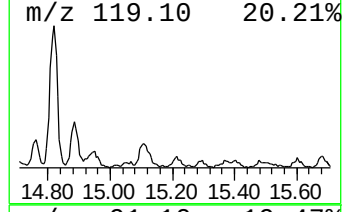
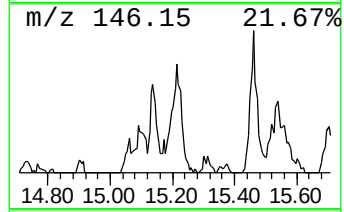
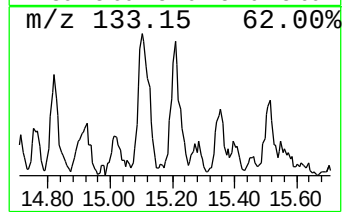
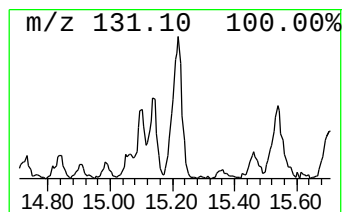
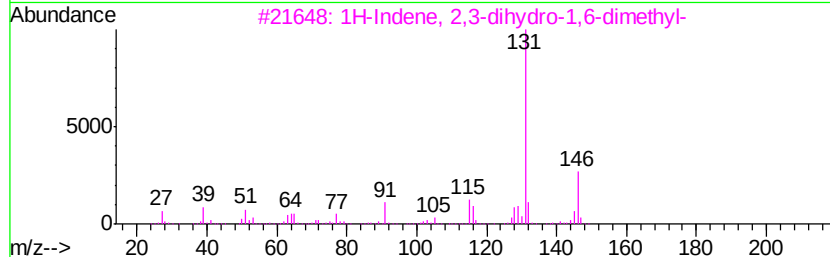
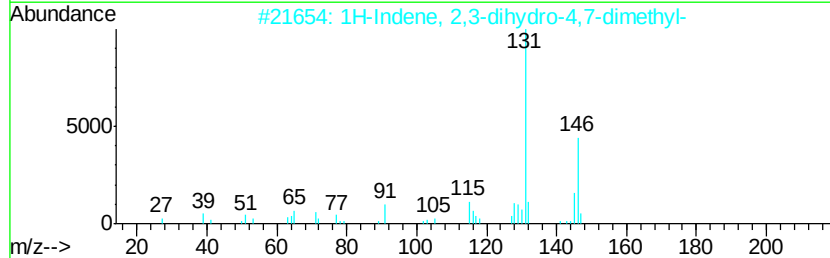
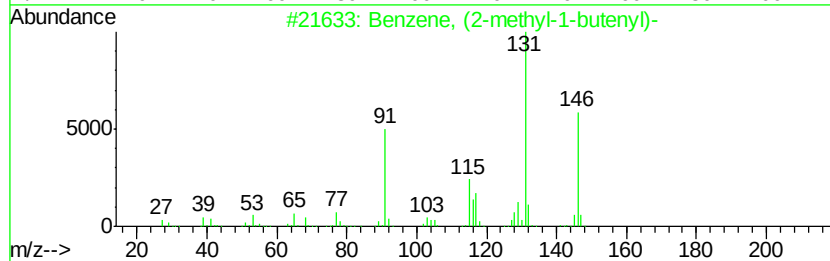
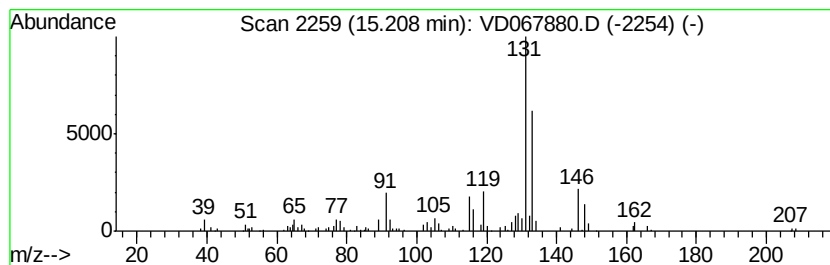
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ClientSampleId :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 5 Benzene, (2-methyl-1-butenyl)- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.21	9.37 ug/l	188603	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6 60
2		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 58
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 53
4		Bicyclo[4.2.0]octa-1,3,5-triene,...	146 C11H14	027087-54-3 49
5		Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1 49



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

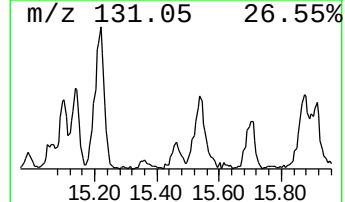
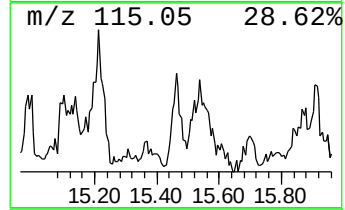
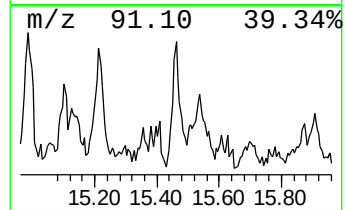
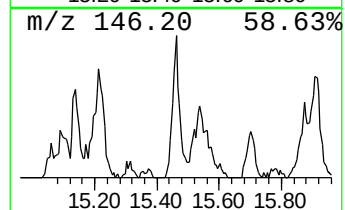
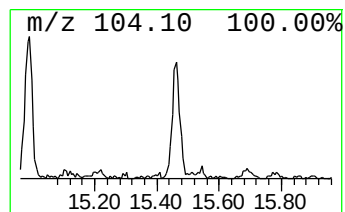
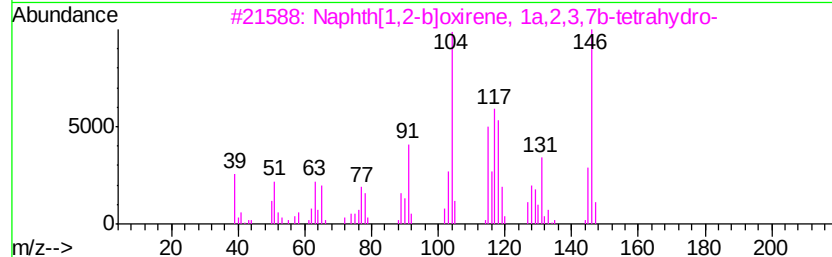
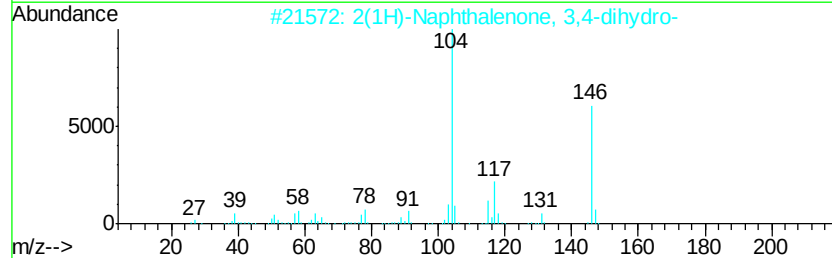
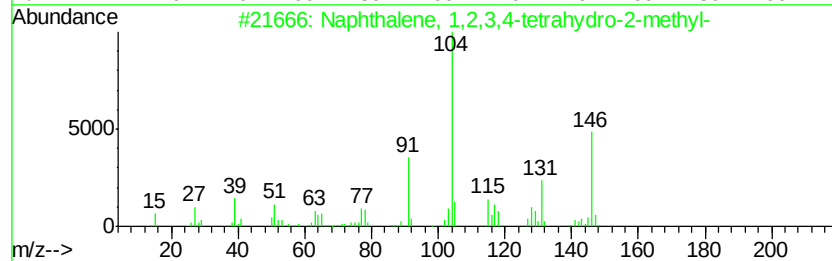
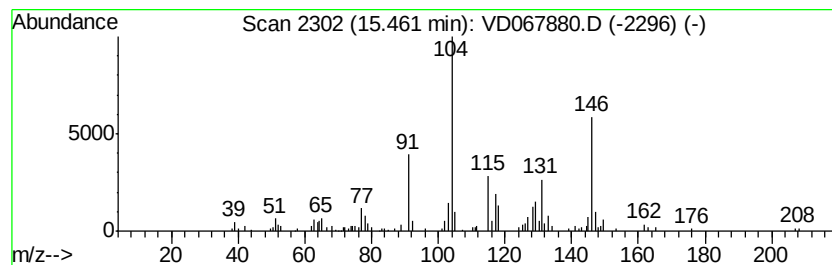
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ClientSampleId :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 6 Naphthalene, 1,2,3,4-tetra... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.46	6.05 ug/l	121631	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	003877-19-8 87
2		2(1H)-Naphthalenone, 3,4-dihydro-	146 C10H10O	000530-93-8 68
3		Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146 C10H10O	002461-34-9 62
4		2-Propyn-1-ol, 3-(4-methylphenyl)-	146 C10H10O	016017-24-6 45
5		7-Azabicyclo[4.2.2]deca-2,4,9-tr...	147 C9H9NO	017198-06-0 43



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

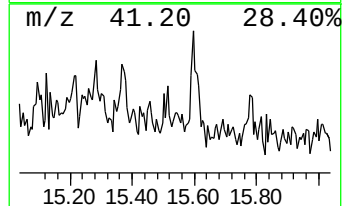
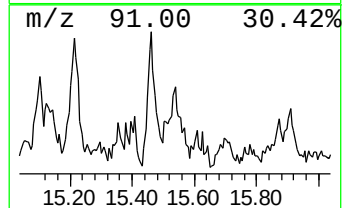
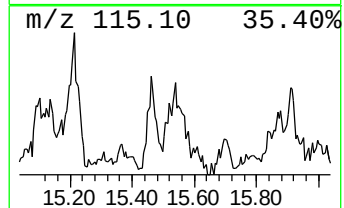
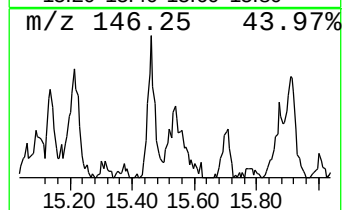
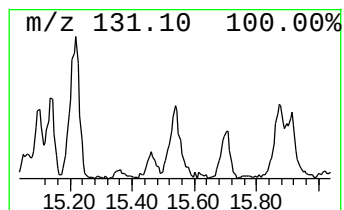
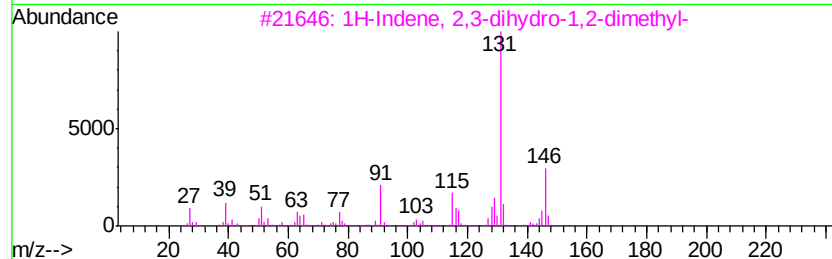
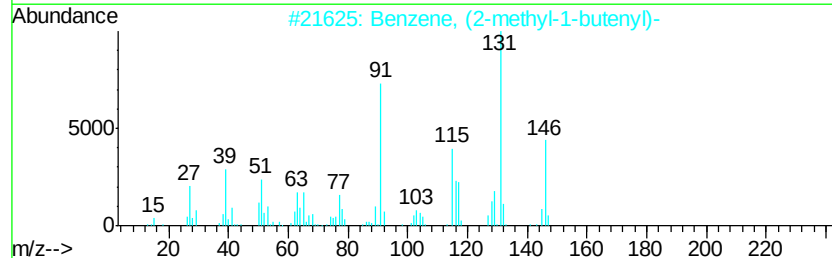
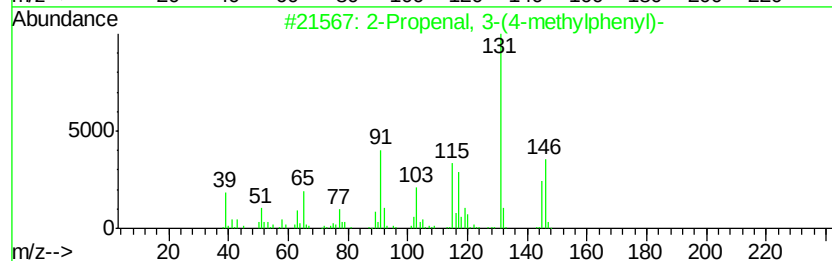
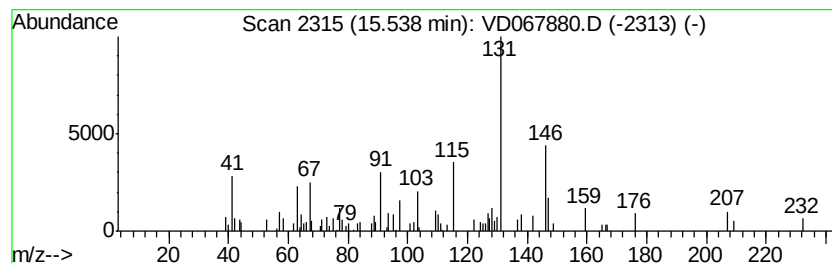
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ClientSampleId :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 7 2-Propenal, 3-(4-methylphen... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.54	6.14 ug/l	123452	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1	2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2	59
2	Benzene, (2-methyl-1-butenyl)-	146 C11H14	056253-64-6	58
3	1H-Indene, 2,3-dihydro-1,2-dimet...	146 C11H14	017057-82-8	53
4	Benzene, 1-methyl-4-(1-methyl-2-...	146 C11H14	097664-18-1	53
5	Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3	50



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-02-12320

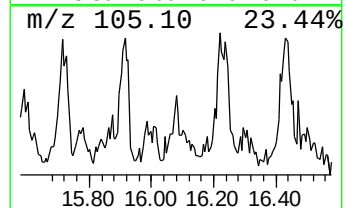
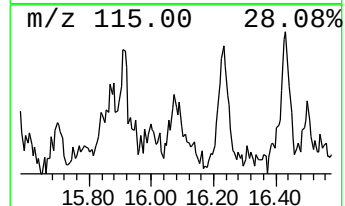
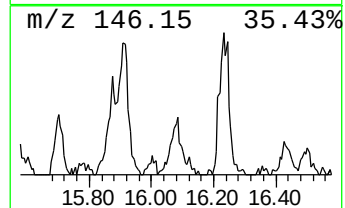
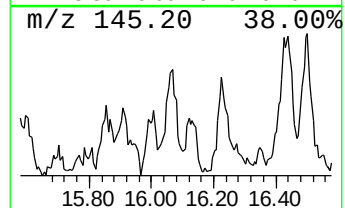
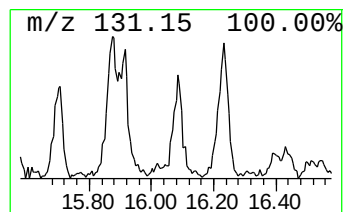
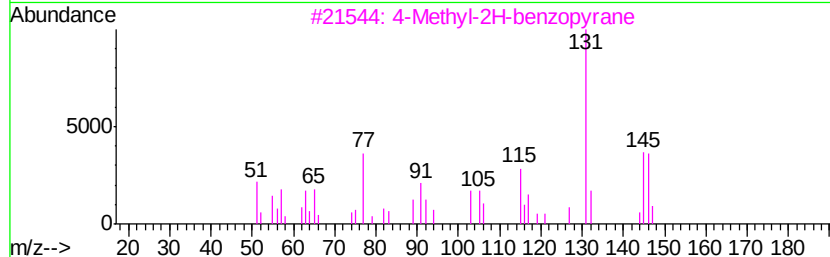
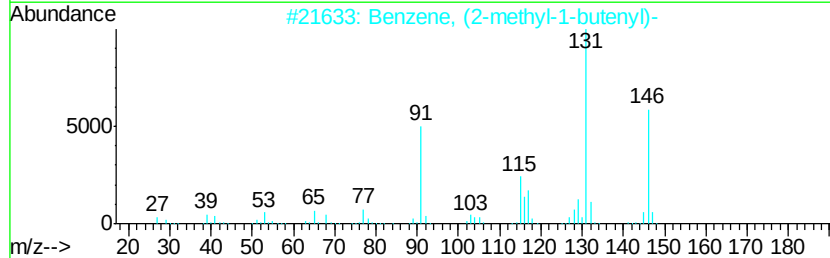
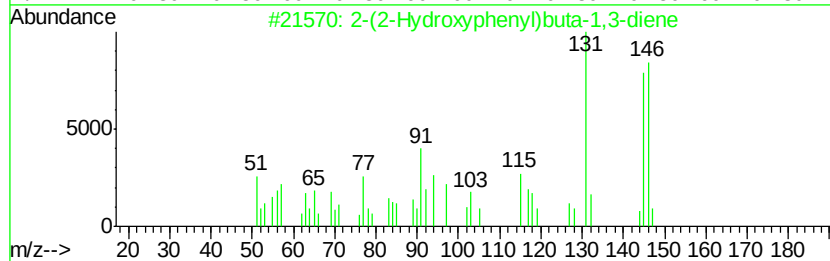
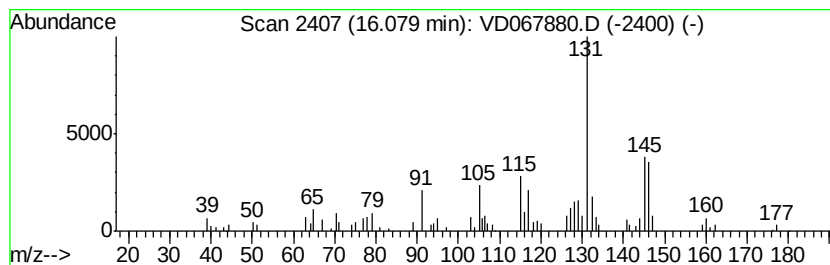
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Quant Title : SW846 8260

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

Peak Number 8 2-(2-Hydroxyphenyl)buta-1,3... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.08	7.34 ug/l	147743	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	2-(2-Hydroxyphenyl)buta-1,3-diene	146	C10H10O	038865-47-3	87
2		Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	83
3		4-Methyl-2H-benzopyrane	146	C10H10O	021776-94-3	81
4		Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	80
5		1,3-Cyclopentadiene, 5-(trans-2-...	146	C11H14	079209-36-2	74



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

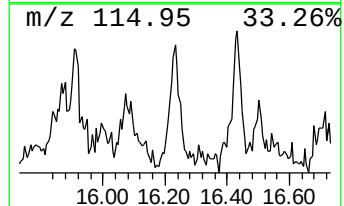
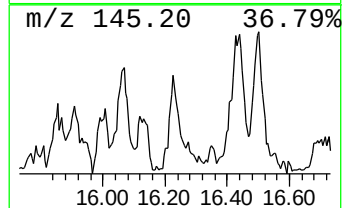
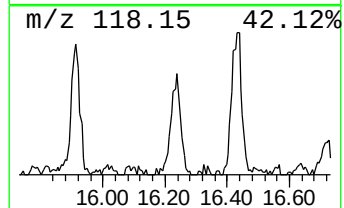
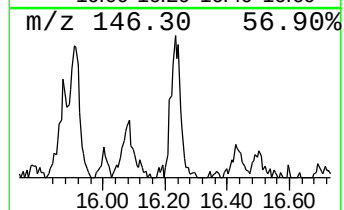
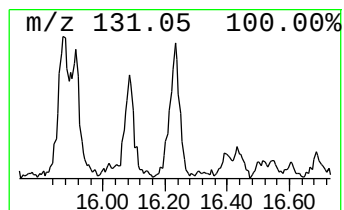
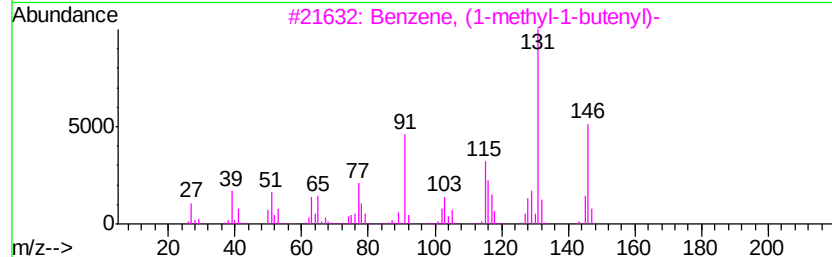
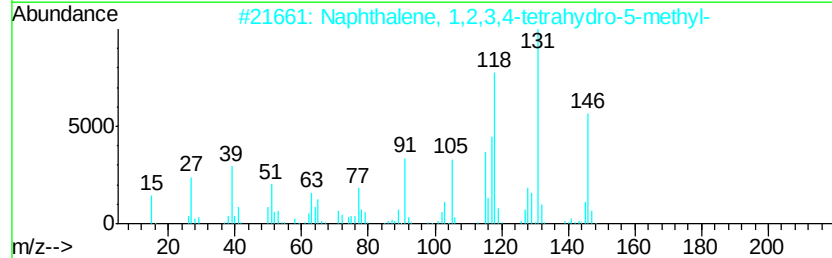
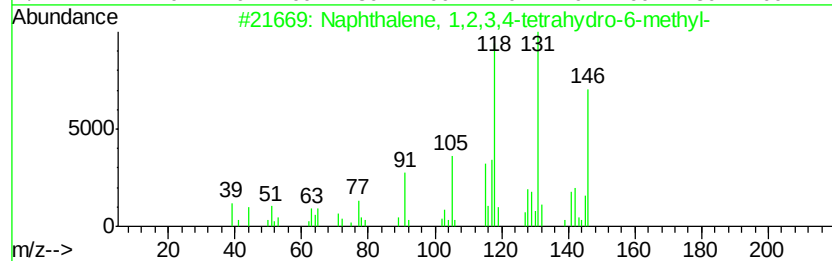
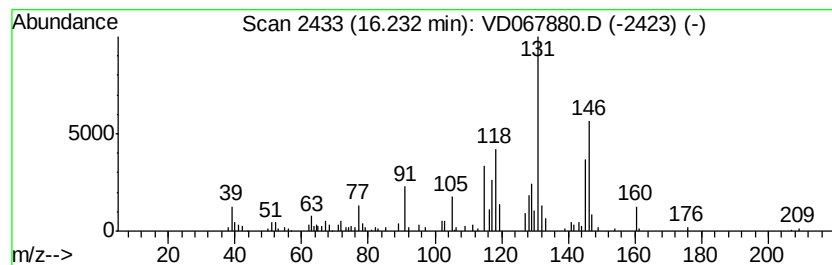
Instrument :
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ClientSampled :
EO-02-12320

Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 9 Naphthalene, 1,2,3,4-tetra... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.23	9.15 ug/l	184174	1,4-Dichlorobenzene-d4	13.58
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001680-51-9 86
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	002809-64-5 83
3		Benzene, (1-methyl-1-butenyl)-	146 C11H14	053172-84-2 76
4		2-Propenal, 3-(4-methylphenyl)-	146 C10H10O	001504-75-2 64
5		Benzene, 2-ethenyl-1,3,5-trimethyl-	146 C11H14	000769-25-5 64



Data Path : Z:\V0ASRV\HPCHEM1\MSVOA D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-02-12320

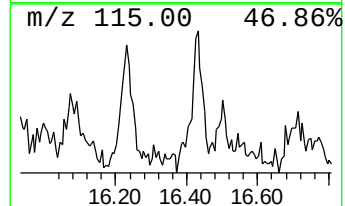
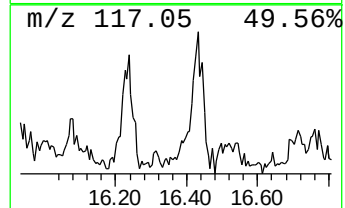
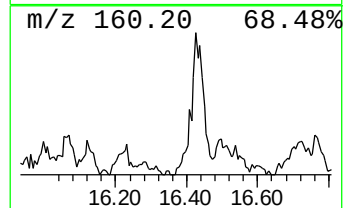
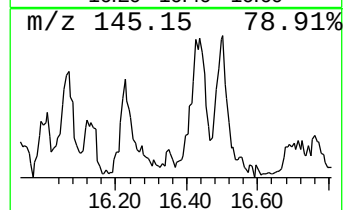
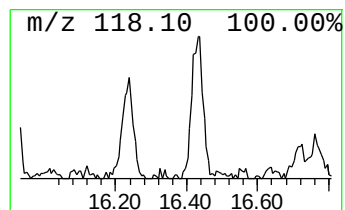
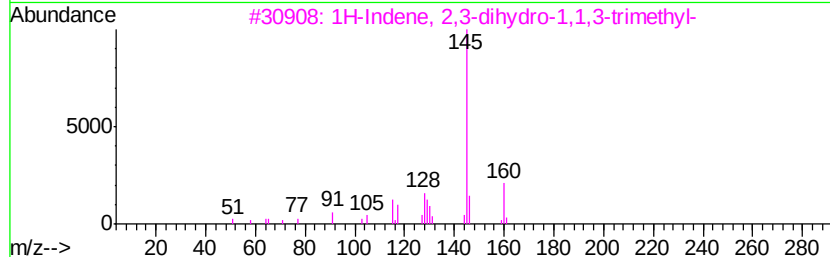
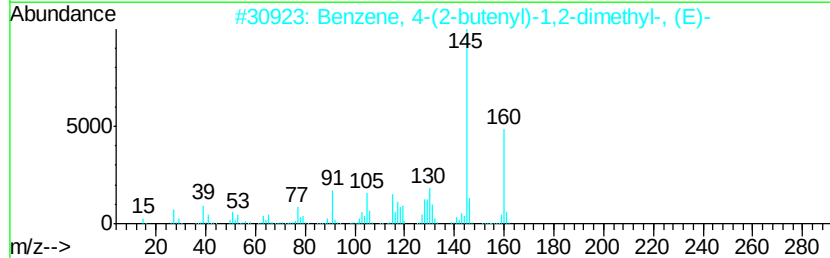
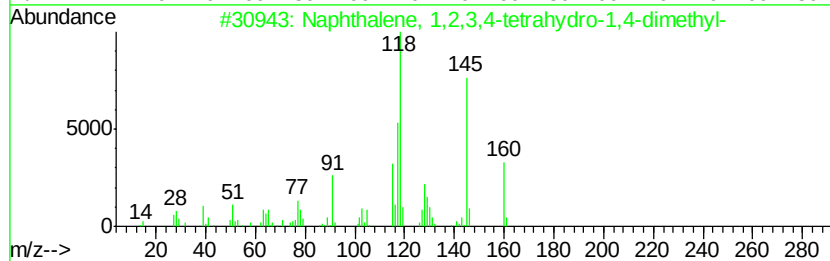
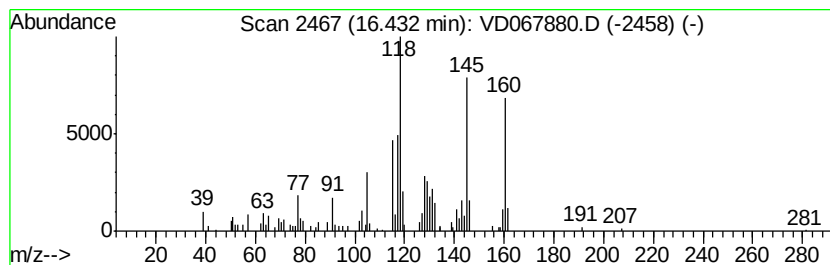
Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.M
Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.43	8.09 ug/l	162721	1,4-Dichlorobenzene-d4	13.58

Hit#	of	Tentative ID	MW	MolForm	CAS#	Qual
1	5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	83
2		Benzene, 4-(2-butenyl)-1,2-dimet...	160	C12H16	054340-86-2	64
3		1H-Indene, 2,3-dihydro-1,1,3-tri...	160	C12H16	002613-76-5	62
4		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021564-91-0	62
5		Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	007524-63-2	58



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD120420\
Data File : VD067880.D
Acq On : 04 Dec 2020 12:30
Operator : VA/SY
Sample : L4939-01
Misc : 7.46G/5.00ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
EO-02-12320

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D120120S.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-ethyl-...	13.81	9.2	ug/l	185456	4	13.58	1005900	50.0
unknown13.90	13.90	12.6	ug/l	254361	4	13.58	1005900	50.0
unknown14.23	14.23	7.8	ug/l	157897	4	13.58	1005900	50.0
o-Cymene	14.82	20.8	ug/l	418680	4	13.58	1005900	50.0
Benzene, (2-methy...	15.21	9.4	ug/l	188603	4	13.58	1005900	50.0
Naphthalene, 1,2,...	15.46	6.0	ug/l	121631	4	13.58	1005900	50.0
2-Propenal, 3-(4-...	15.54	6.1	ug/l	123452	4	13.58	1005900	50.0
2-(2-Hydroxypheny...	16.08	7.3	ug/l	147743	4	13.58	1005900	50.0
Naphthalene, 1,2,...	16.23	9.2	ug/l	184174	4	13.58	1005900	50.0
Naphthalene, 1,2,...	16.43	8.1	ug/l	162721	4	13.58	1005900	50.0