

Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
 Data File : VD067977.D
 Acq On : 30 Dec 2020 15:38
 Operator : VA/SY
 Sample : L5244-01
 Misc : 5.81G/5.00ml/MSVOA_D/SOIL
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_D
 ClientSampleId :
 HR-01-123020

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\82D122220S.M

Title : Sw846 8260

Signal : TIC: VD067977.D\data.ms

peak #	R.T. min	first scan	max scan	last scan	PK TY	peak height	corr. area	corr. % max.	% of total
1	4.956	508	516	525	rVB6	6258	16824	1.23%	0.095%
2	7.920	1009	1020	1025	rBV2	197632	499334	36.36%	2.824%
3	7.979	1025	1030	1041	rVB	279278	631632	45.99%	3.572%
4	8.332	1079	1090	1102	rBV	186034	432251	31.47%	2.444%
5	8.867	1172	1181	1191	rBV	480502	962042	70.05%	5.440%
6	10.343	1423	1432	1439	rBV	766682	1373344	100.00%	7.766%
7	10.402	1439	1442	1449	rVB4	12613	23080	1.68%	0.131%
8	11.643	1643	1653	1666	rBV	650188	1120013	81.55%	6.334%
9	11.855	1679	1689	1692	rBV3	23434	48623	3.54%	0.275%
10	12.179	1732	1744	1753	rBV4	42100	110292	8.03%	0.624%
11	12.355	1770	1774	1776	rBV4	10246	14573	1.06%	0.082%
12	12.390	1776	1780	1787	rVB8	16410	35822	2.61%	0.203%
13	12.632	1815	1821	1831	rVB	473424	775546	56.47%	4.386%
14	12.796	1845	1849	1856	rVB6	22809	49227	3.58%	0.278%
15	12.885	1856	1864	1866	rBV2	66541	116136	8.46%	0.657%
16	12.914	1866	1869	1872	rBV	42460	51142	3.72%	0.289%
17	12.961	1872	1877	1883	rVB3	187872	322883	23.51%	1.826%
18	13.008	1883	1885	1890	rVB4	24863	32780	2.39%	0.185%
19	13.120	1890	1904	1910	rBV	116452	252760	18.40%	1.429%
20	13.185	1911	1915	1923	rVB5	39337	77888	5.67%	0.440%
21	13.267	1923	1929	1934	rBV	112353	179275	13.05%	1.014%
22	13.396	1948	1951	1955	rVB5	15970	20275	1.48%	0.115%
23	13.461	1955	1962	1967	rBV3	110061	217963	15.87%	1.233%
24	13.508	1967	1970	1975	rVB3	39248	60458	4.40%	0.342%
25	13.573	1975	1981	1985	rBV	553111	1042176	75.89%	5.893%
26	13.773	2010	2015	2018	rBV2	196280	280158	20.40%	1.584%
27	13.808	2018	2021	2031	rVB2	293236	529195	38.53%	2.993%
28	13.896	2031	2036	2042	rBV3	161594	289166	21.06%	1.635%
29	13.967	2042	2048	2055	rBV	160474	269000	19.59%	1.521%
30	14.032	2055	2059	2061	rBV3	71133	101118	7.36%	0.572%
31	14.061	2061	2064	2069	rVB2	142698	196018	14.27%	1.108%
32	14.120	2069	2074	2078	rVB	266409	419802	30.57%	2.374%
33	14.179	2078	2084	2087	rBV3	35232	58974	4.29%	0.333%
34	14.226	2087	2092	2097	rVB3	279357	455775	33.19%	2.577%
35	14.284	2097	2102	2103	rBV2	37854	60452	4.40%	0.342%
36	14.361	2109	2115	2118	rBV2	129851	218170	15.89%	1.234%
37	14.408	2119	2123	2125	rVV3	101747	153888	11.21%	0.870%

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Instrument :
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 ClientSampleId :
 HR-01-123020

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\82D122220S.M
 Title : Sw846 8260

38	14.443	2125	2129	2131	rVV	141532	224100	16.32%	1.267%
39	14.479	2131	2135	2139	rVV	263399	385596	28.08%	2.181%
40	14.520	2139	2142	2146	rVV	133675	186783	13.60%	1.056%
41	14.561	2146	2149	2157	rVV5	119486	203314	14.80%	1.150%
42	14.661	2157	2166	2171	rVV2	123980	247574	18.03%	1.400%
43	14.708	2171	2174	2179	rVV5	97753	188295	13.71%	1.065%
44	14.755	2179	2182	2187	rVB	128918	189118	13.77%	1.069%
45	14.820	2187	2193	2200	rBV5	462806	1186880	86.42%	6.712%
46	14.884	2200	2204	2209	rVB2	86785	141637	10.31%	0.801%
47	14.984	2215	2221	2227	rVB	271284	454076	33.06%	2.568%
48	15.055	2229	2233	2235	rBV5	26137	41428	3.02%	0.234%
49	15.096	2235	2240	2243	rBV4	130603	225650	16.43%	1.276%
50	15.208	2252	2259	2266	rVB3	178697	438883	31.96%	2.482%
51	15.296	2271	2274	2280	rVB5	25975	45800	3.33%	0.259%
52	15.367	2280	2286	2291	rBV9	35965	93008	6.77%	0.526%
53	15.455	2295	2301	2306	rBV3	145834	271656	19.78%	1.536%
54	15.508	2306	2310	2311	rBV4	55364	67314	4.90%	0.381%
55	15.590	2323	2324	2333	rVB5	51723	91964	6.70%	0.520%
56	15.702	2334	2343	2349	rBV6	62139	167576	12.20%	0.948%
57	15.773	2351	2355	2360	rVV7	21985	45760	3.33%	0.259%
58	15.867	2363	2371	2372	rVV4	57008	121126	8.82%	0.685%
59	15.908	2373	2378	2384	rVB	169841	318505	23.19%	1.801%
60	16.002	2390	2394	2398	rVV6	18973	30066	2.19%	0.170%
61	16.067	2399	2405	2421	rVB5	52509	202325	14.73%	1.144%
62	16.226	2423	2432	2441	rBV2	126184	290383	21.14%	1.642%
63	16.426	2456	2466	2472	rBV4	80064	194982	14.20%	1.103%
64	16.490	2473	2477	2483	rVB4	27498	49684	3.62%	0.281%
65	16.702	2504	2513	2518	rBV4	29254	82190	5.98%	0.465%

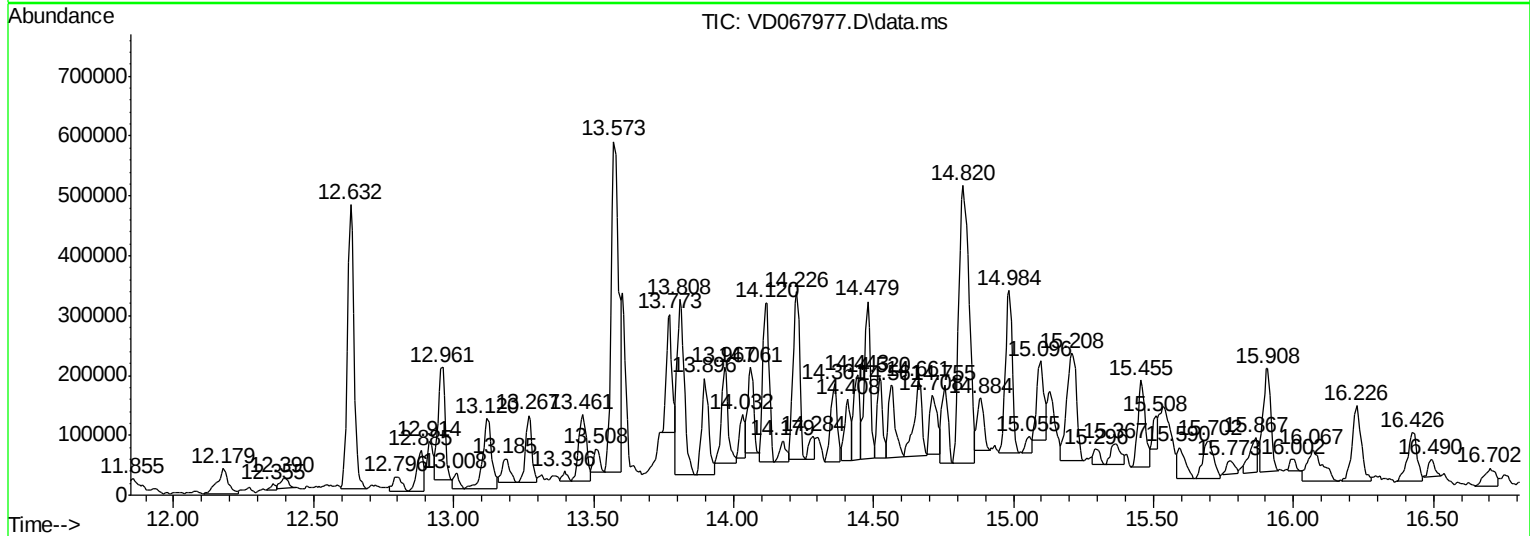
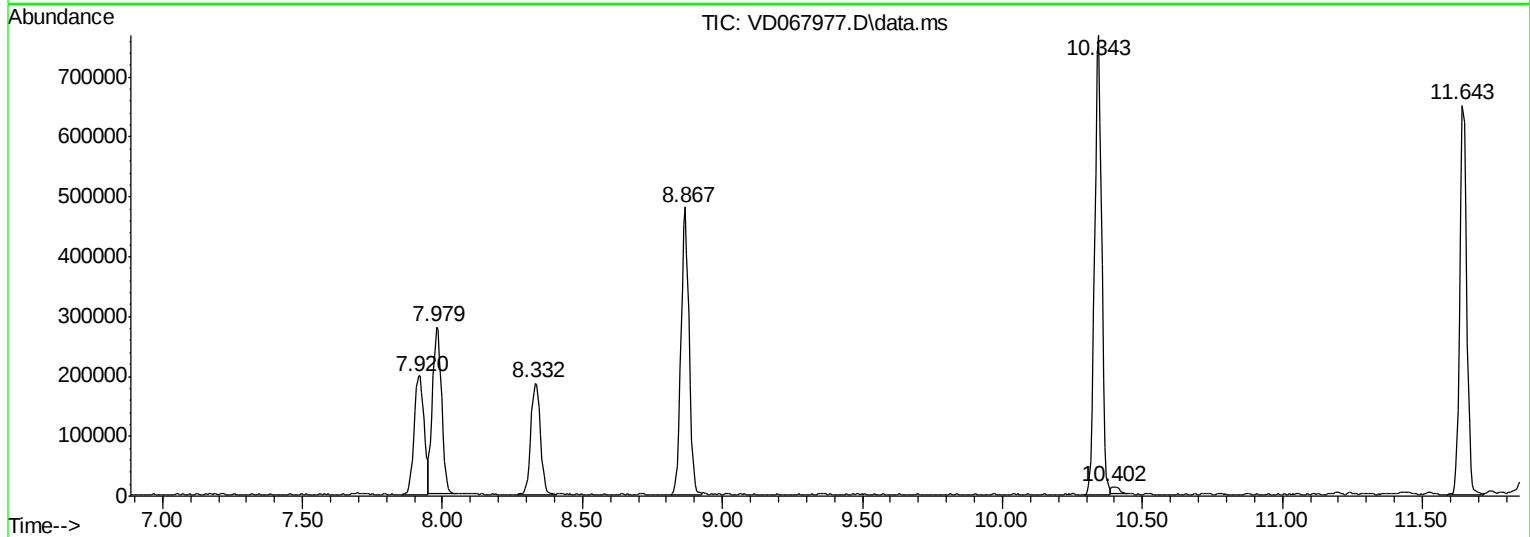
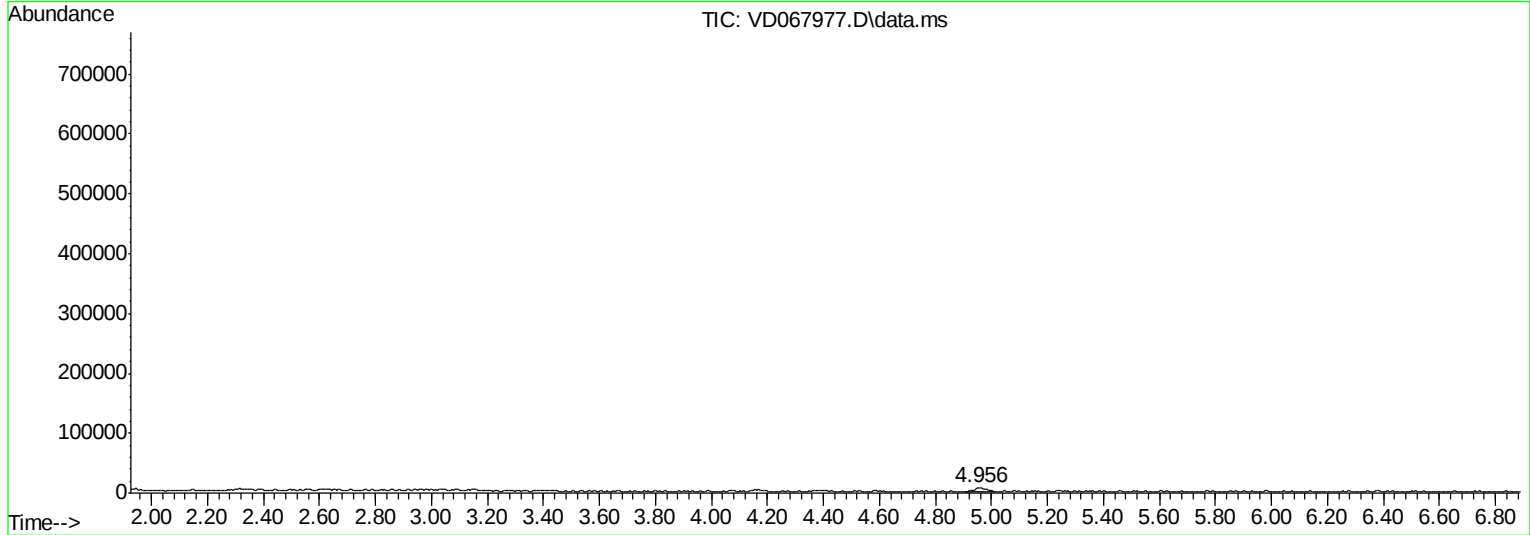
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

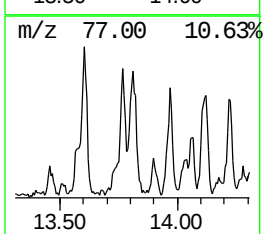
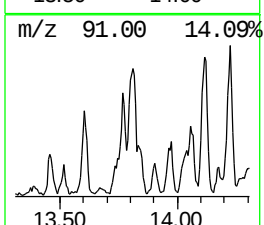
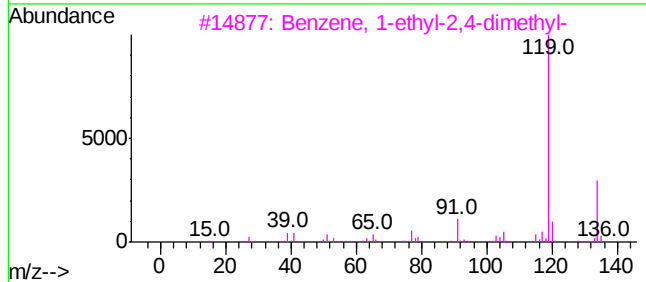
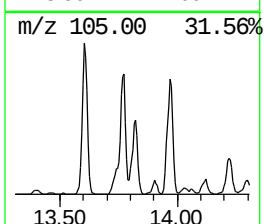
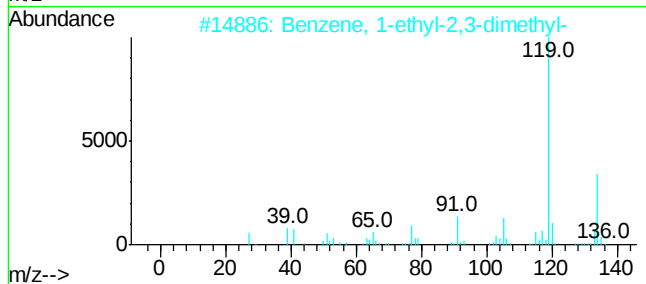
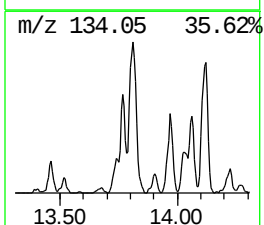
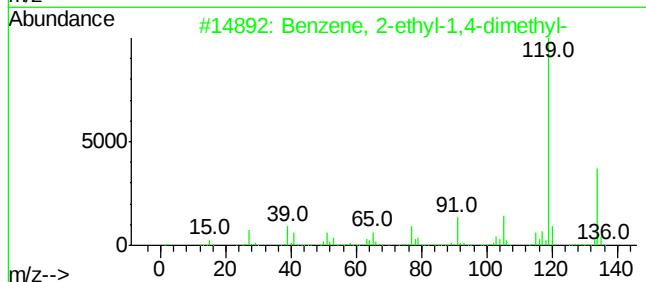
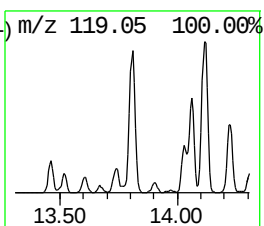
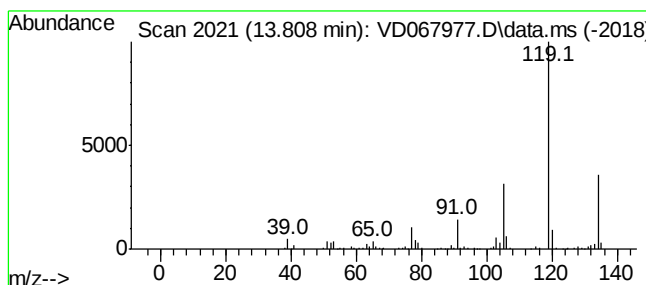
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.808	25.39 ug/l	529195	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	87
2	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	86
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	83
4	Benzene, 1-ethyl-3,5-dimethyl-	134	C10H14	000934-74-7	80
5	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	80



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Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

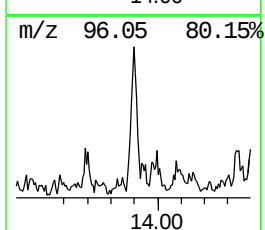
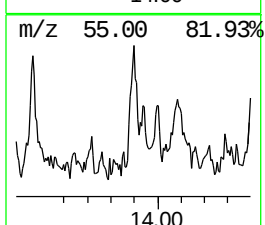
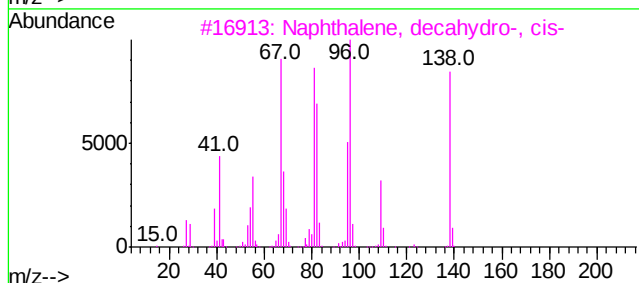
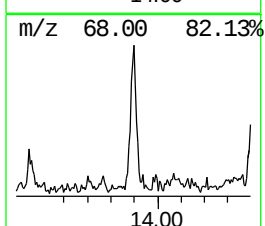
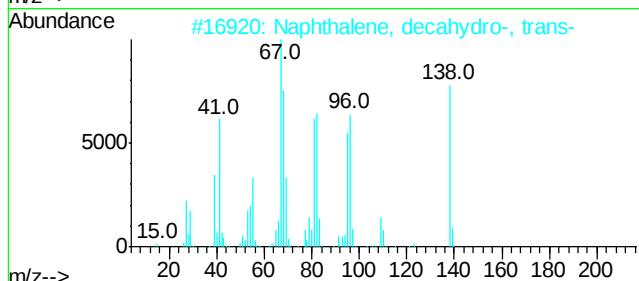
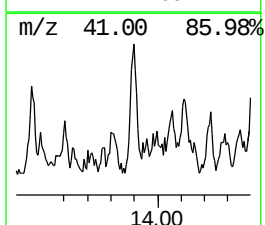
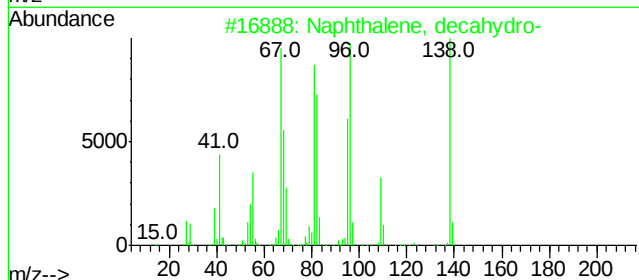
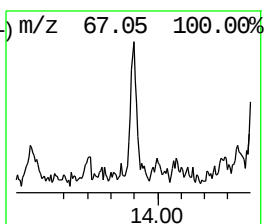
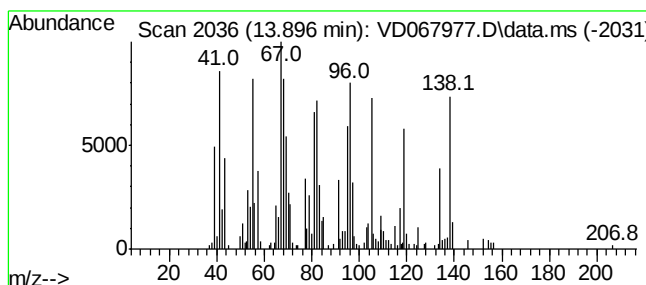
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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

Peak Number 2 Naphthalene, decahydro- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.896	13.87 ug/l	289166	1,4-Dichlorobenzene-d4	13.573

Hit#	of	5	Tentative ID	Mw	MolForm	CAS#	Qual
1			Naphthalene, decahydro-	138	C10H18	000091-17-8	96
2			Naphthalene, decahydro-, trans-	138	C10H18	000493-02-7	95
3			Naphthalene, decahydro-, cis-	138	C10H18	000493-01-6	92
4			Spiro[4.5]decane	138	C10H18	000176-63-6	64
5			Cyclohexane, 1-methyl-4-(1-methy...	138	C10H18	001124-25-0	55



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

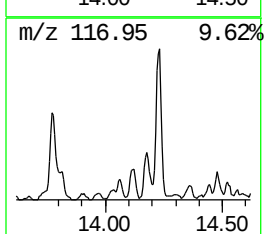
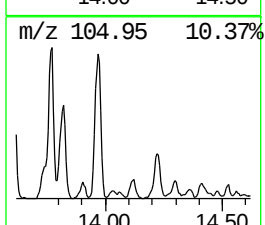
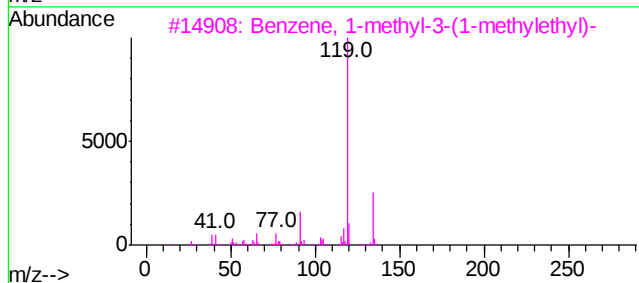
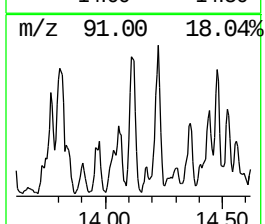
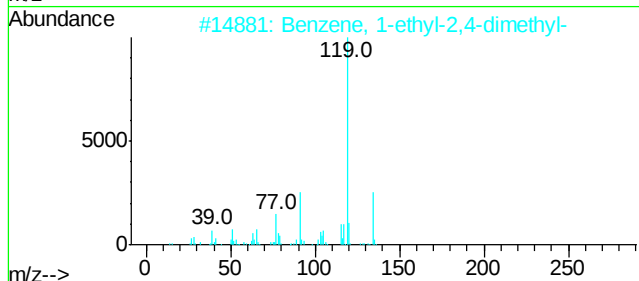
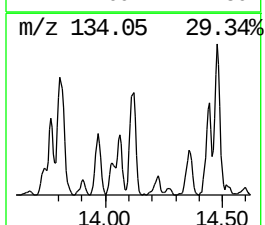
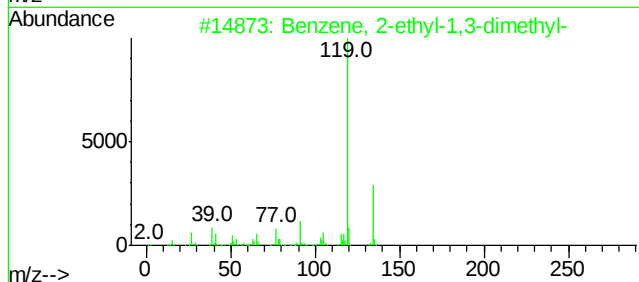
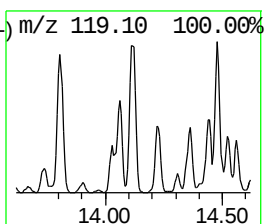
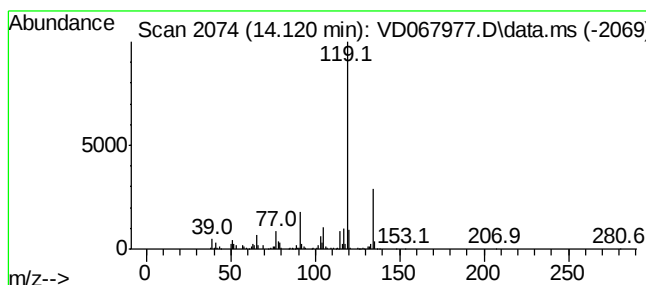
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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

Peak Number 3 Benzene, 2-ethyl-1,3-dimethyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.120	20.14 ug/l	419802	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
2	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	p-Cymene	134	C10H14	000099-87-6	95



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ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

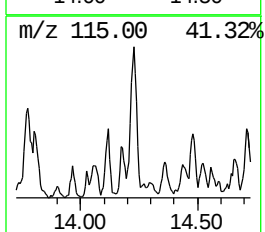
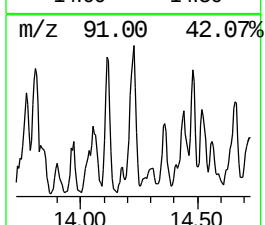
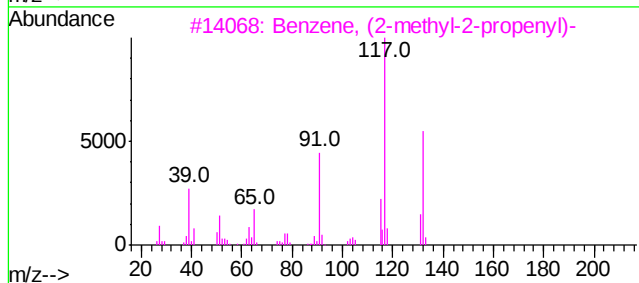
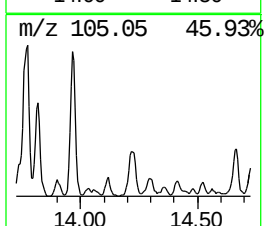
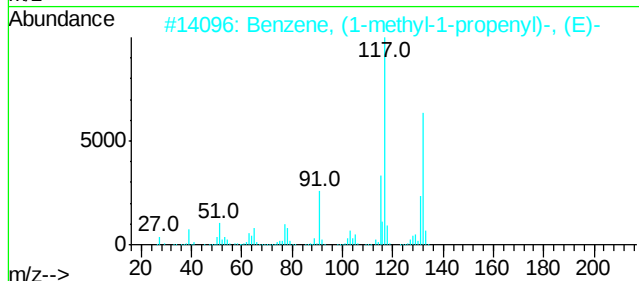
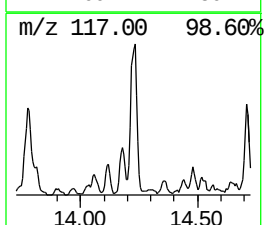
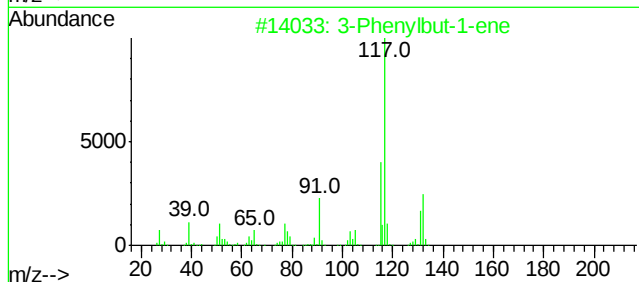
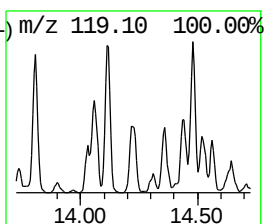
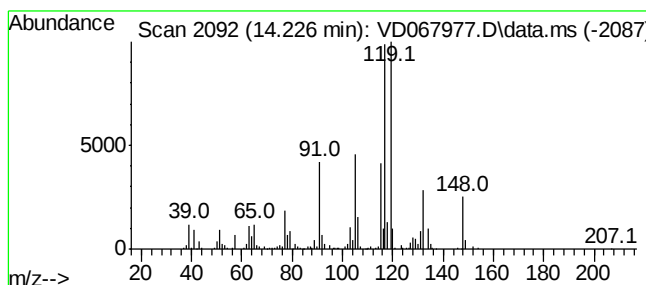
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.M
Quant Title : SW846 8260

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

Peak Number 4 3-Phenylbut-1-ene Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.226	21.87 ug/l	455775	1,4-Dichlorobenzene-d4	13.573

Hit# of	5	Tentative ID	Mw	MolForm	CAS#	Qual
1	3-Phenylbut-1-ene	132	C10H12	000934-10-1	55	
2	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	55	
3	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	46	
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000767-99-7	46	
5	3a,6-Methano-3aH-indene, 2,3,6,7...	132	C10H12	098640-29-0	46	



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
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Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-01-123020

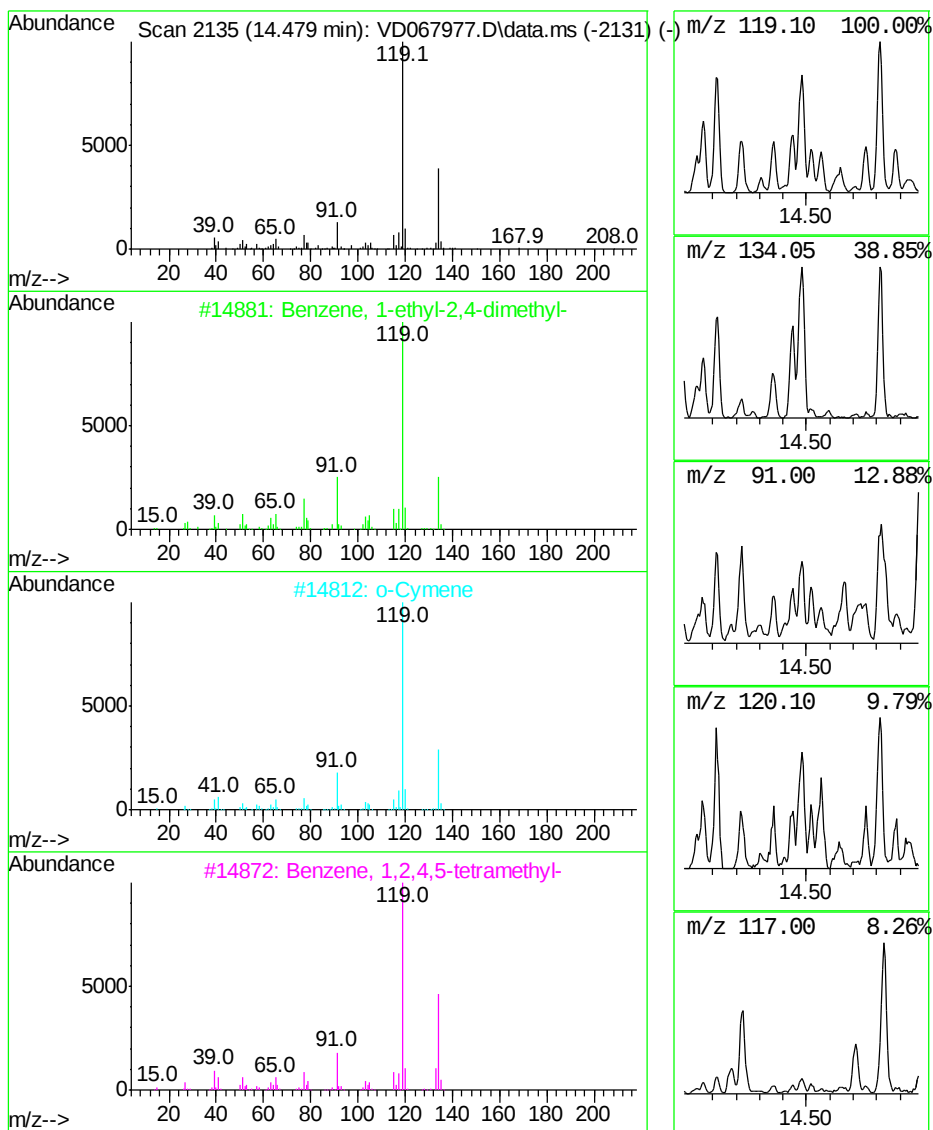
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TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

Peak Number 5 Benzene, 1-ethyl-2,4-dimethyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.479	18.50 ug/l	385596	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	94
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	94
4	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	91
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	91



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
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ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-01-123020

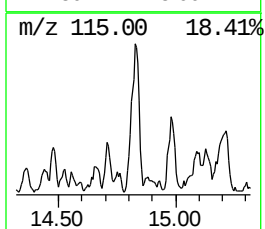
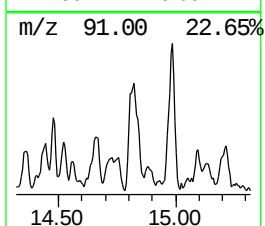
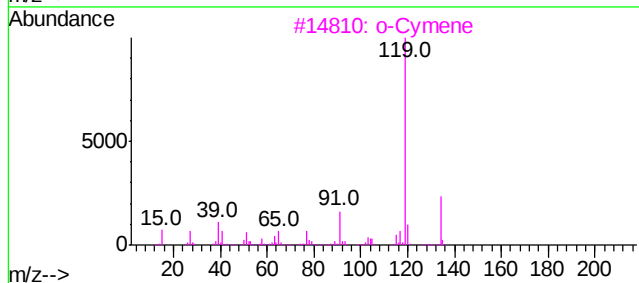
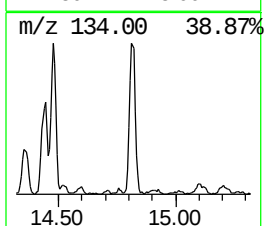
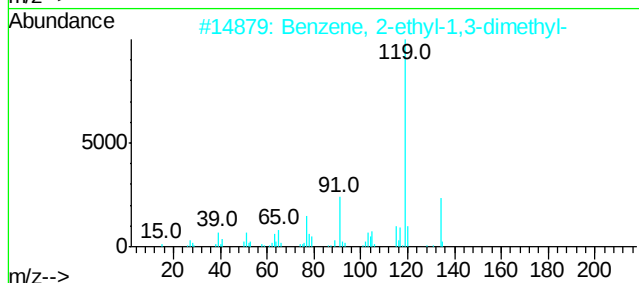
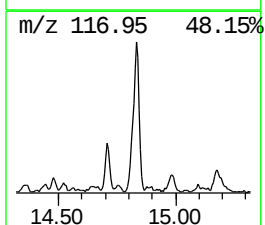
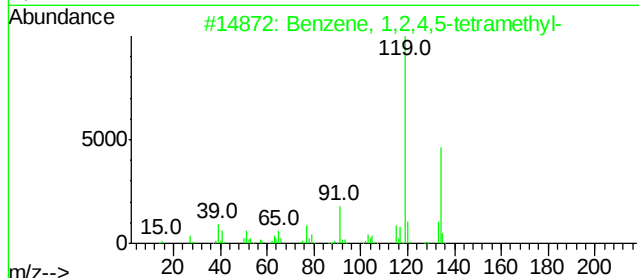
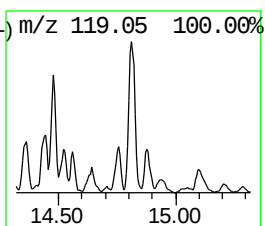
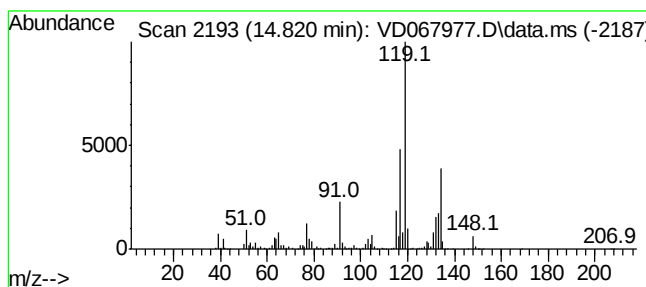
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Quant Title : SW846 8260

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

Peak Number 6 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.820	56.94 ug/l	1186880	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	70
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	64
3	o-Cymene	134	C10H14	000527-84-4	64
4	p-Cymene	134	C10H14	000099-87-6	64
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	64



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

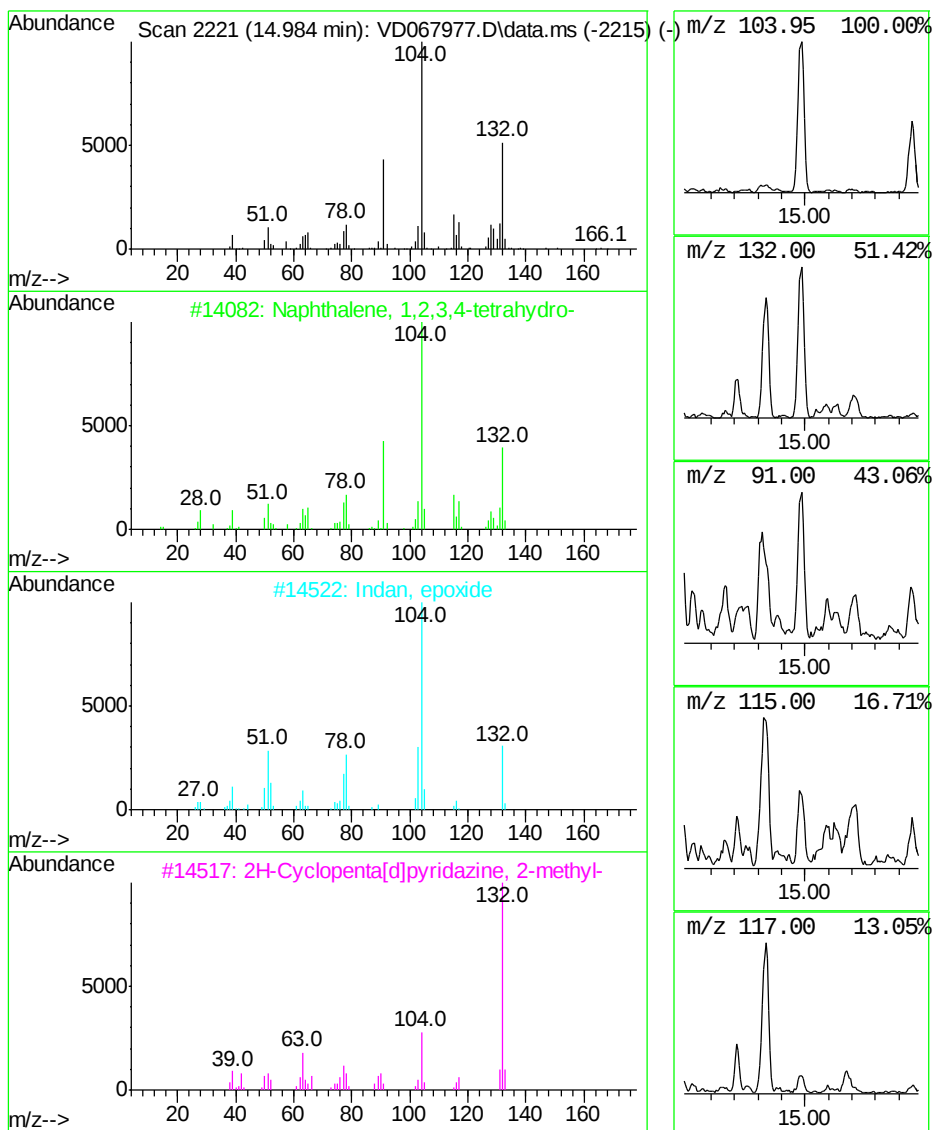
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.984	21.79 ug/l	454076	1,4-Dichlorobenzene-d4	13.573

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-Cyclopenta[d]pyridazine, 2-me...	132	C8H8N2	022291-85-6	47
4	2H-Inden-2-one, 1,3-dihydro-	132	C9H8O	000615-13-4	35
5	Benzene, 1,1'-(1,2-cyclobutanedi...	208	C16H16	007694-30-6	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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ClientSampleId :
HR-01-123020

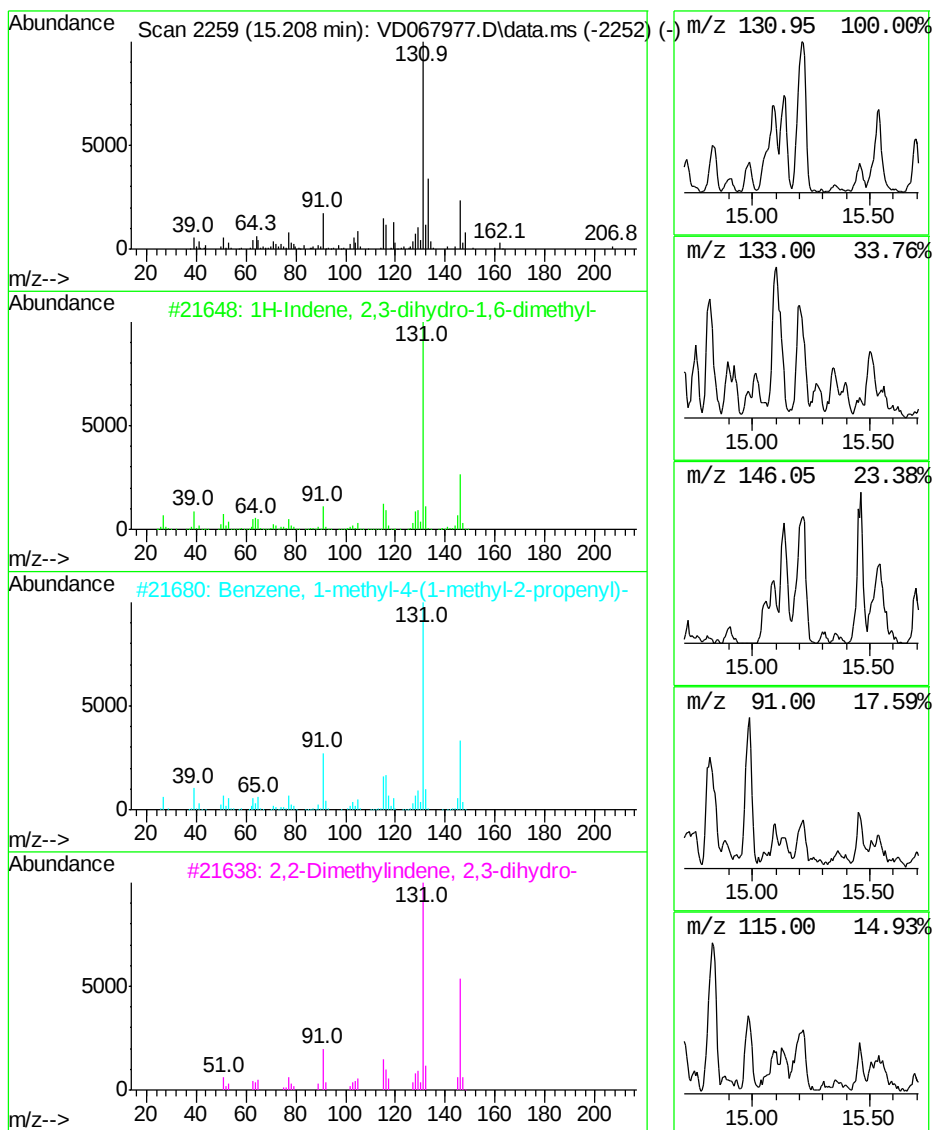
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Quant Title : SW846 8260

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

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.208	21.06 ug/l	438883	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	89
2	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	81
3	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	76
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
5	Benzene, (1-methyl-1-butenyl)-	146	C11H14	053172-84-2	76



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
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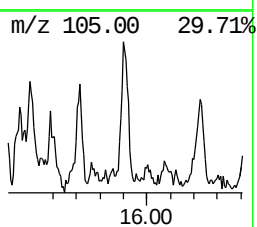
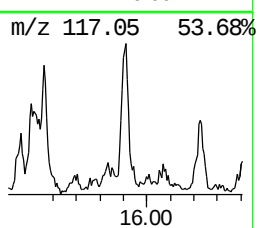
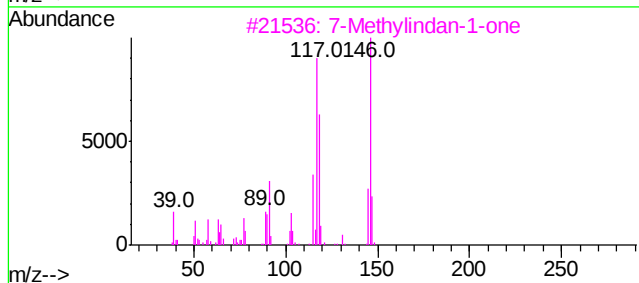
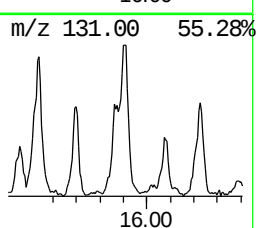
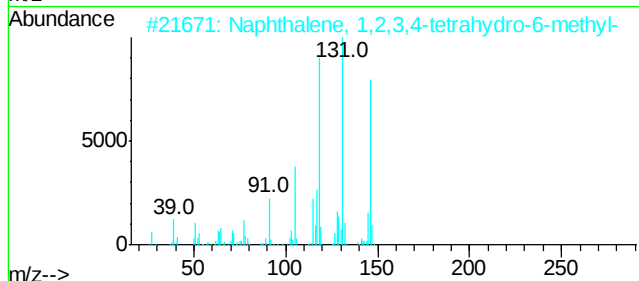
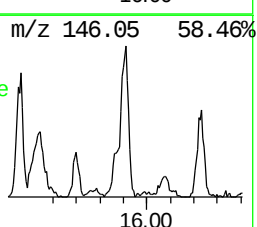
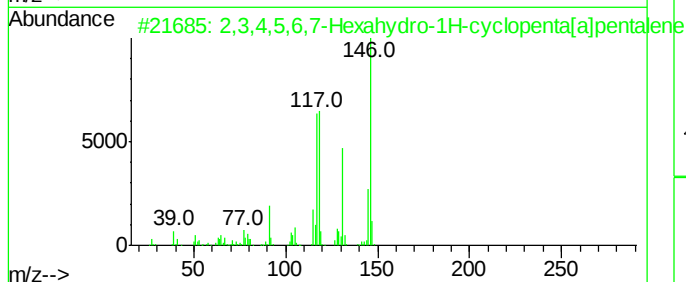
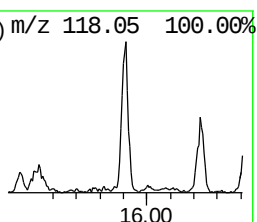
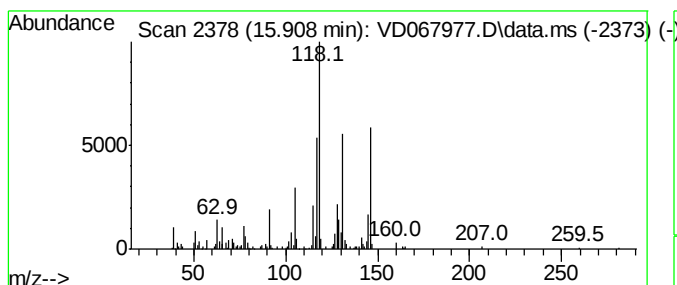
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Quant Title : SW846 8260

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

Peak Number 9 2,3,4,5,6,7-Hexahydro-1H-cy... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.908	15.28 ug/l	318505	1,4-Dichlorobenzene-d4	13.573

Hit# of 5	Tentative ID	Mw	MolForm	CAS#	Qual
1	2,3,4,5,6,7-Hexahydro-1H-cyclope...	146	C11H14	1000189-31-0	53
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	47
3	7-Methylindan-1-one	146	C10H10O	039627-61-7	41
4	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	38
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	35



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

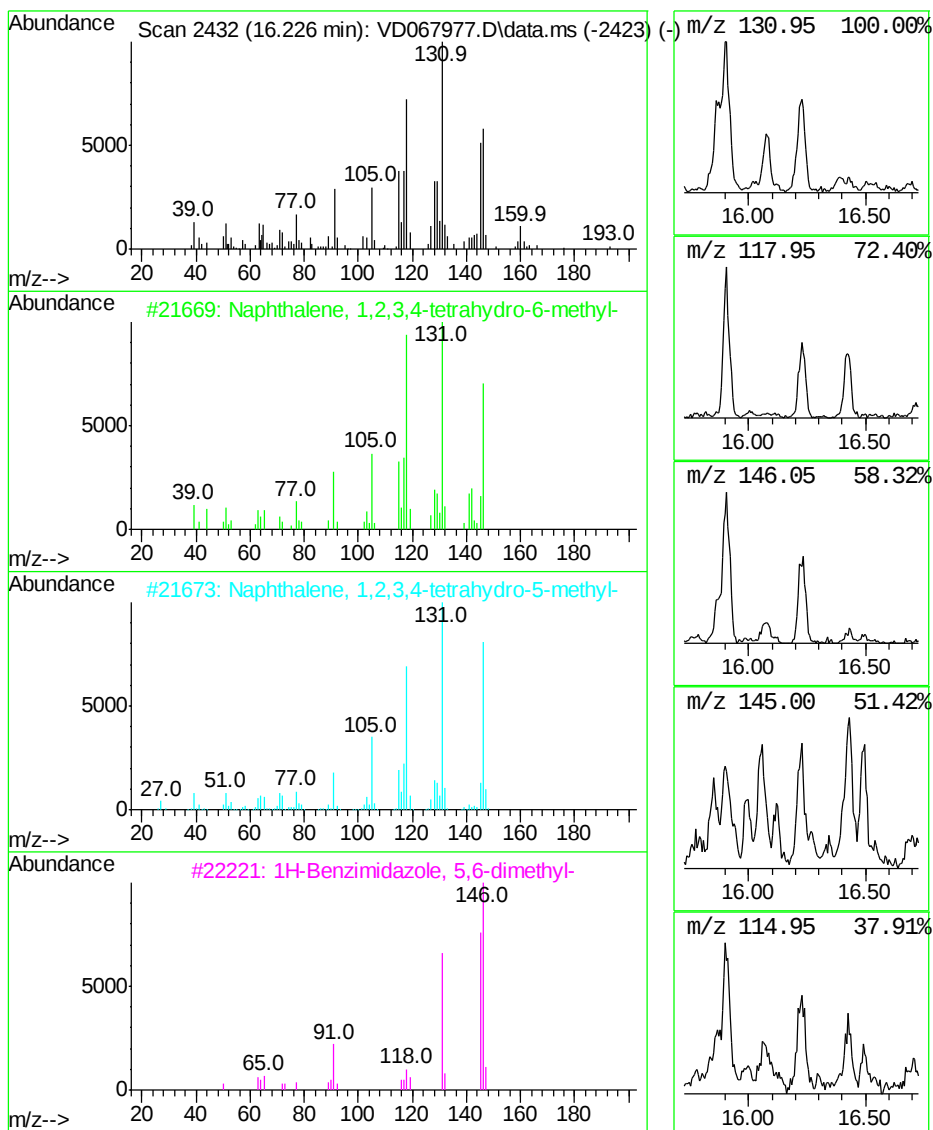
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Quant Title : SW846 8260

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

Peak Number 10 Naphthalene, 1,2,3,4-tetrahydro... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.226	13.93 ug/l	290383	1,4-Dichlorobenzene-d4	13.573

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	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	60
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	60
5	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	53



Data Path : Z:\voasrv\HPCHEM1\MSVOA_D\Data\VD123020\
Data File : VD067977.D
Acq On : 30 Dec 2020 15:38
Operator : VA/SY
Sample : L5244-01
Misc : 5.81G/5.00ml/MSVOA_D/SOIL
ALS Vial : 9 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
HR-01-123020

Quant Method : Z:\voasrv\HPCHEM1\MSVOA_D\Method\82D122220S.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-ethy...	13.808	25.4	ug/l	529195	4	13.573	1042180	50.0
Naphthalene, de...	13.896	13.9	ug/l	289166	4	13.573	1042180	50.0
Benzene, 2-ethy...	14.120	20.1	ug/l	419802	4	13.573	1042180	50.0
3-Phenylbut-1-ene	14.226	21.9	ug/l	455775	4	13.573	1042180	50.0
Benzene, 1-ethy...	14.479	18.5	ug/l	385596	4	13.573	1042180	50.0
Benzene, 1,2,4,...	14.820	56.9	ug/l	1186880	4	13.573	1042180	50.0
Naphthalene, 1,...	14.984	21.8	ug/l	454076	4	13.573	1042180	50.0
1H-Indene, 2,3-...	15.208	21.1	ug/l	438883	4	13.573	1042180	50.0
2,3,4,5,6,7-Hex...	15.908	15.3	ug/l	318505	4	13.573	1042180	50.0
Naphthalene, 1,...	16.226	13.9	ug/l	290383	4	13.573	1042180	50.0