LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

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\82D110618S.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	1.715	17	19	23	rVB	54067	86537	1.99%	0.198%
2	1.852	30	33	44	rVB2	68917	146035	3.36%	0.333%
3	2.344	79	83	90	rBV2	25268	57642	1.33%	0.132%
4	3.151	159	165	176	rVB3	17281	68219	1.57%	0.156%
5	3.722	216	223	238	rBV2	248264	812397	18.72%	1.854%
6	4.106	256	262	272	rBV2	28455	91891	2.12%	0.210%
7	4.519	298	304	312	rBV	42296	118815	2.74%	0.271%
8	5.405	390	394	401	rVB	37422	102459	2.36%	0.234%
9	5.651	407	419	428	rBV2	28840	102264	2.36%	0.233%
10	6.104	460	465	472	rVB5	16584	44061	1.02%	0.101%
11	6.330	479	488	492	rBV	678189	1883441	43.40%	4.299%
12	6.409	492	496	512	rVB	1406361	3844609	88.58%	8.776%
13	6.615	512	517	521	rBV	26612	72636	1.67%	0.166%
14	6.792	529	535	544	rBV	453405	1140474	26.28%	2.603%
15	6.930	544	549	551	rVV3	25845	75070	1.73%	0.171%
16	6.989	551	555	561	rVV2	50224	173452	4.00%	0.396%
17	7.088	561	565	573	rVB2	27763	77924	1.80%	0.178%
18	7.442	595	601	618	rBV	1813055	4340198	100.00%	9.907%
19	8.042	653	662	670	rBV2	130586	434919	10.02%	0.993%
20	8.150	670	673	678	rVB2	20222	50544	1.16%	0.115%
21	8.436	693	702	706	rBV4	15357	51764	1.19%	0.118%
22	8.643	718	723	730	rVB3	39740	109071	2.51%	0.249%
23	8.829	732	742	749	rBV2	40847	200109	4.61%	0.457%
24	9.213	774	781	785	rVB3	15706	46392	1.07%	0.106%
25	9.459	800	806	821	rVB	1800680	4094076	94.33%	9.345%
26	9.951	850	856	861	rVB4	30338	81195	1.87%	0.185%
27	10.109	866	872	877	rBV3	44416	112741	2.60%	0.257%
28	10.188	877	880	884	rBV2	20480	45974	1.06%	0.105%
29	10.699	929	932	936	rVB3	27997	56280	1.30%	0.128%
30	10.778	936	940	944	rBV2	34719	72768	1.68%	0.166%
31	11.044	963	967	970	rBV2	26613	50428	1.16%	0.115%
32	11.280	987	991	998	rVV	2547734	4291518	98.88%	9.796%
33	11.555	1013	1019	1021	rBV6	22577	59041	1.36%	0.135%
34	11.890	1048	1053	1056	rBV4	25078	62075	1.43%	0.142%

LSC Area Percent Report

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Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

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\82D110618S.M
Title : SW846 8260

35	11.929	1056	1057	1066	rVB5	24431	59346	1.37%	0.135%
36	12.106	1066	1075	1077	rBV4	14124	47553	1.10%	0.109%
37	12.205	1079	1085	1087	rVV4	17645	47842	1.10%	0.109%
38	12.431	1104	1108	1116	rBV	1736354	2596655	59.83%	5.927%
39	12.569	1116	1122	1126	rBV2	44571	109817	2.53%	0.251%
40	12.795	1142	1145	1151	rVB4	46544	117485	2.71%	0.268%
41	12.943	1156	1160	1162	rBV	113891	160714	3.70%	0.367%
42	13.219	1182	1188	1193	rBV4	28656	73544	1.69%	0.168%
43	13.297	1193	1196	1200	rVV5	55511	138820	3.20%	0.317%
44	13.376	1200	1204	1206	rVV	2630661	3650070	84.10%	8.332%
45	13.405	1206	1207	1212	rVV	210522	318641	7.34%	0.727%
46	13.484	1212	1215	1219	rVV4	44477	108955	2.51%	0.249%
47	13.553	1219	1222	1227	rVV2	146858	302589	6.97%	0.691%
48	13.622	1227	1229	1231	rVV	211935	305865	7.05%	0.698%
49	13.661	1231	1233	1235	rVV3	113955	191736	4.42%	0.438%
50	13.701	1235	1237	1239	rVV	87775	133820	3.08%	0.305%
51	13.760	1240	1243	1248	rVV	177306	328953	7.58%	0.751%
52	13.858	1251	1253	1255	rVV3	57778	110049	2.54%	0.251%
53	13.898	1255	1257	1260	rVV	345606	534363	12.31%	1.220%
54	13.937	1260	1261	1264	rVV3	86592	151258	3.49%	0.345%
55	13.986	1264	1266	1271	rVV2	274733	578785	13.34%	1.321%
56	14.075	1271	1275	1277	rVV3	113720	260866	6.01%	0.595%
57	14.114	1277	1279	1283	rVV2	181800	436446	10.06%	0.996%
58	14.193	1285	1287	1289	rVV	228797	329042	7.58%	0.751%
59	14.222	1289	1290	1293	rVV	139826	181648	4.19%	0.415%
60	14.271	1293	1295	1298	rVV2	243874	451974	10.41%	1.032%
61	14.311	1298	1299	1303	rVV2	141581	290621	6.70%	0.663%
62	14.380	1303	1306	1309	rVV4	129034	360620	8.31%	0.823%
63	14.429	1309	1311	1314	rVV3	113590	266056	6.13%	0.607%
64	14.478	1314	1316	1317	rVV	141512	205190	4.73%	0.468%
65	14.518	1317	1320	1325	rVV2	601430	1157601	26.67%	2.642%
66	14.577	1325	1326	1329	rVV2	179852	290682	6.70%	0.664%
67	14.645	1329	1333	1336	rVV3	233427	492207	11.34%	1.123%
68	14.714	1338	1340	1341	rVV	98342	151172	3.48%	0.345%
69	14.754	1341	1344	1345	rVV2	258093	449565	10.36%	1.026%
70	14.773	1345	1346	1348	rVV	294620	346514	7.98%	0.791%
71	14.842	1348	1353	1359	rVV2	322271	1003449	23.12%	2.290%

LSC Area Percent Report

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ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

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

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Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Title : SW846 8260

72	15.019	1362	1371	1373	rVV5	266139	795606	18.33%	1.816%
73	15.069	1373	1376	1378	rVV	244773	489425	11.28%	1.117%
74	15.108	1378	1380	1382	rVV2	149299	265351	6.11%	0.606%
75	15.137	1382	1383	1388	rVV4	104586	231462	5.33%	0.528%
76	15.206	1388	1390	1393	rVV2	130690	256825	5.92%	0.586%
77	15.265	1393	1396	1398	rVV2	76348	170292	3.92%	0.389%
78	15.324	1398	1402	1403	rVV2	184667	335785	7.74%	0.766%
79	15.354	1403	1405	1409	rVV2	158146	288587	6.65%	0.659%
80	15.423	1409	1412	1415	rVV3	87802	194508	4.48%	0.444%
81	15.472	1415	1417	1420	rVV	138091	245269	5.65%	0.560%
82	15.511	1420	1421	1424	rVV2	57848	73343	1.69%	0.167%
83	15.580	1424	1428	1431	rVV	154898	252438	5.82%	0.576%
84	15.620	1431	1432	1437	rVV5	24402	60092	1.38%	0.137%
85	15.728	1438	1443	1445	rVV3	102219	202127	4.66%	0.461%
86	15.767	1445	1447	1450	rVB	39426	59111	1.36%	0.135%
87	15.905	1458	1461	1465	rBV5	21181	48407	1.12%	0.110%
88	16.013	1469	1472	1477	rVB5	21454	44017	1.01%	0.100%

Sum of corrected areas: 43810177

LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
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Acq On : 19 Nov 2018 14:34
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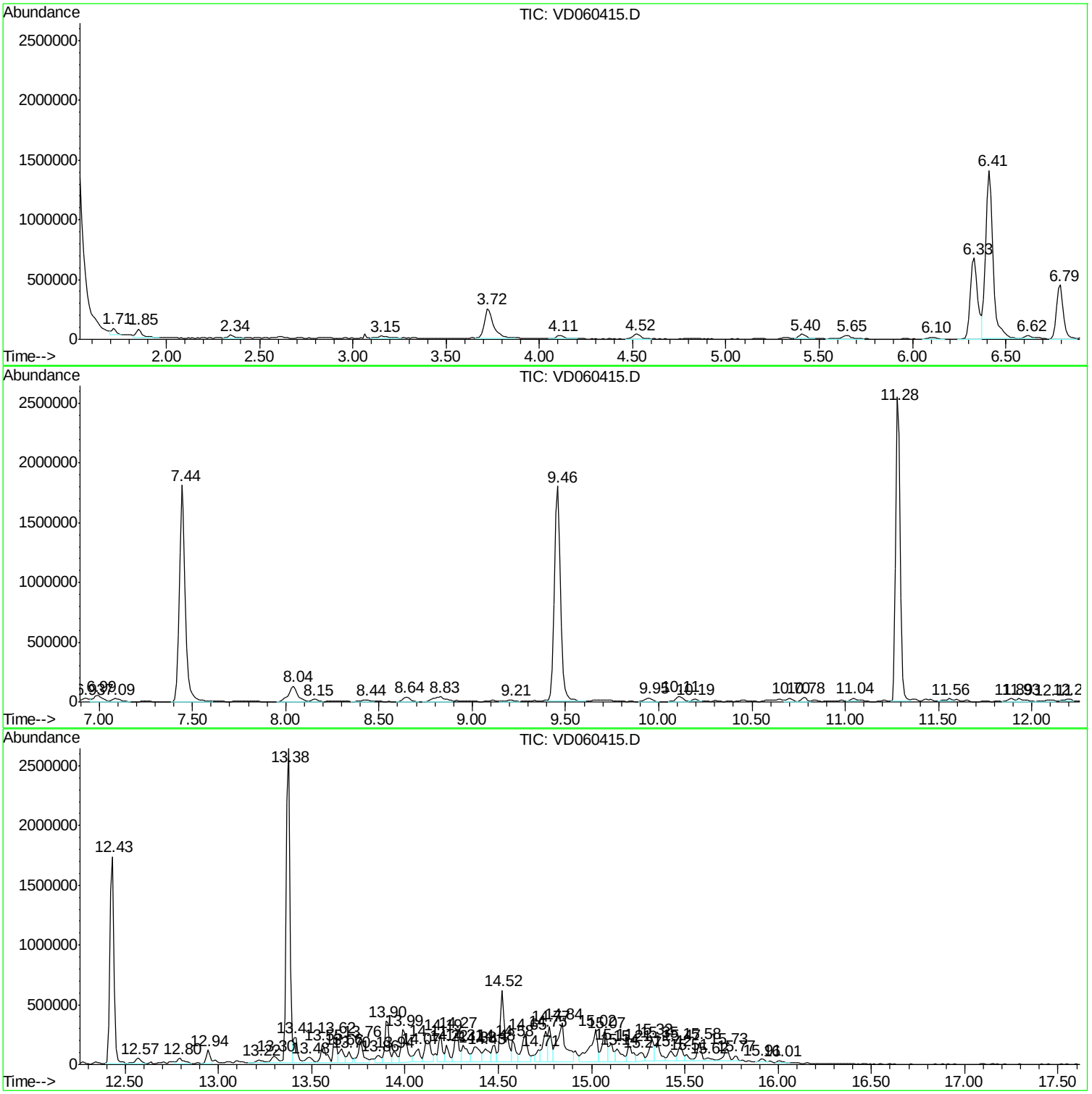
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument :
MSVOA_D
ClientSampleId :
SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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



Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

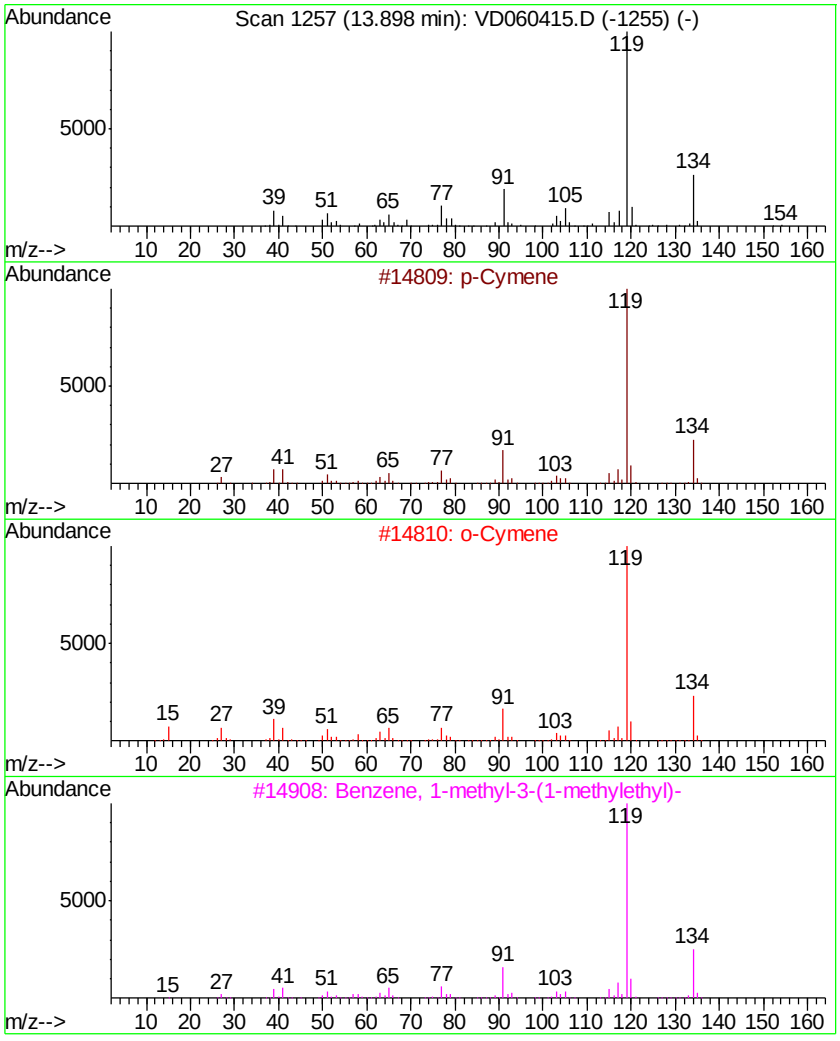
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

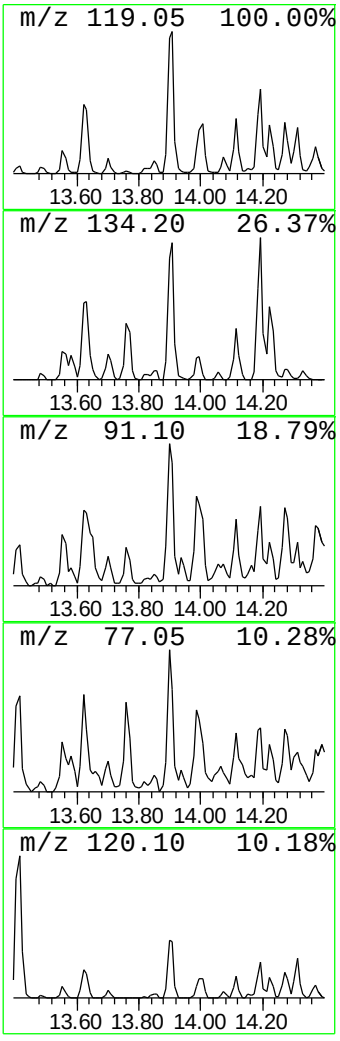
****

Peak Number 1 p-Cymene Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.90	7.32 ug/l	534363	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	p-Cymene	134	C10H14	000099-87-6	94
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	92
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	91





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

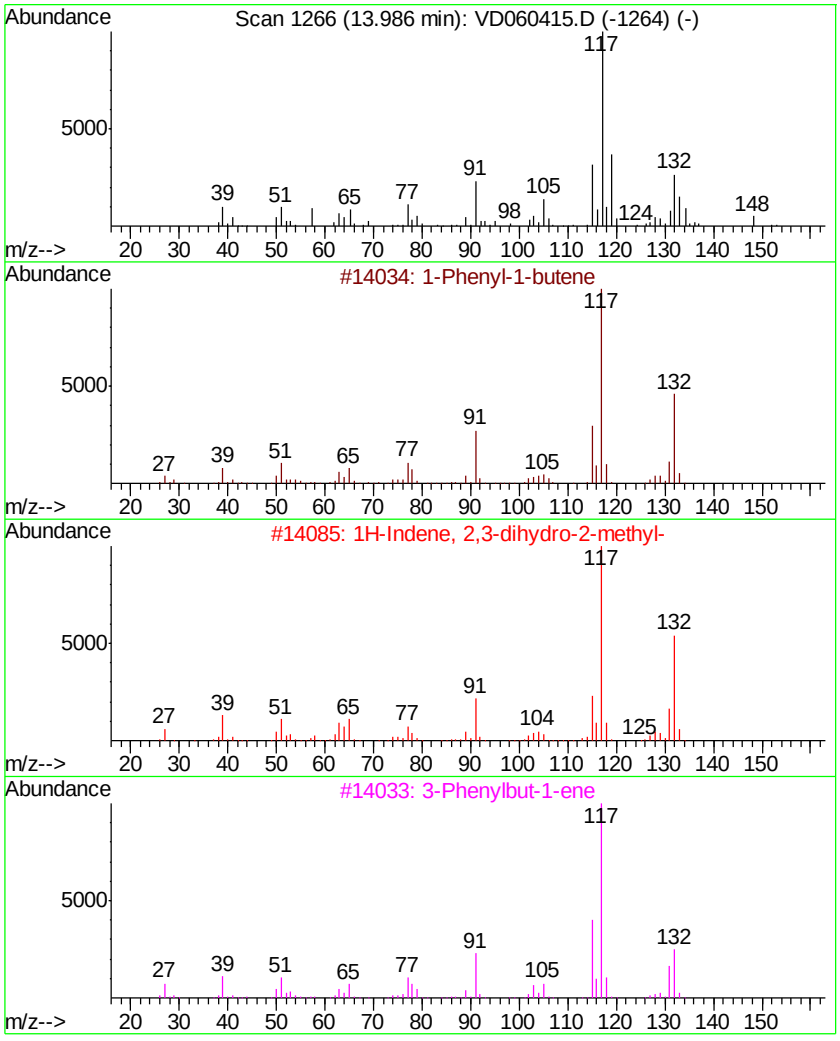
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

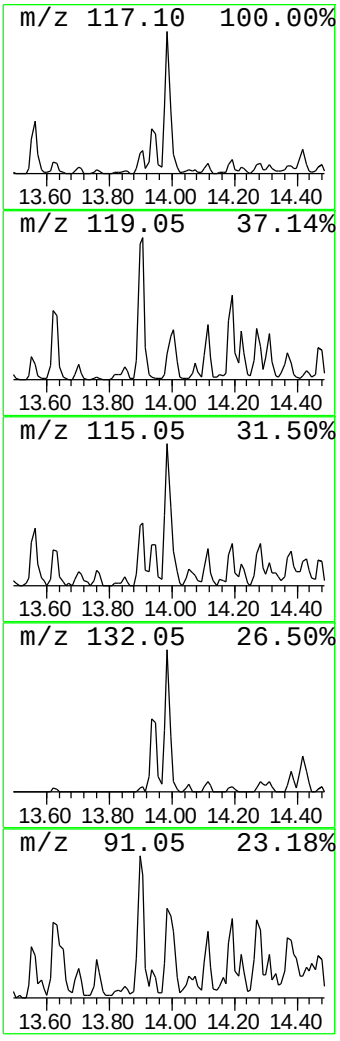
****

Peak Number 2 1-Phenyl-1-butene Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	7.93 ug/l	578785	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	89
2	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	76
3	3-Phenylbut-1-ene	132	C10H12	000934-10-1	76
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
5	Benzene, 1-ethenyl-4-ethyl-	132	C10H12	003454-07-7	76





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
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ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

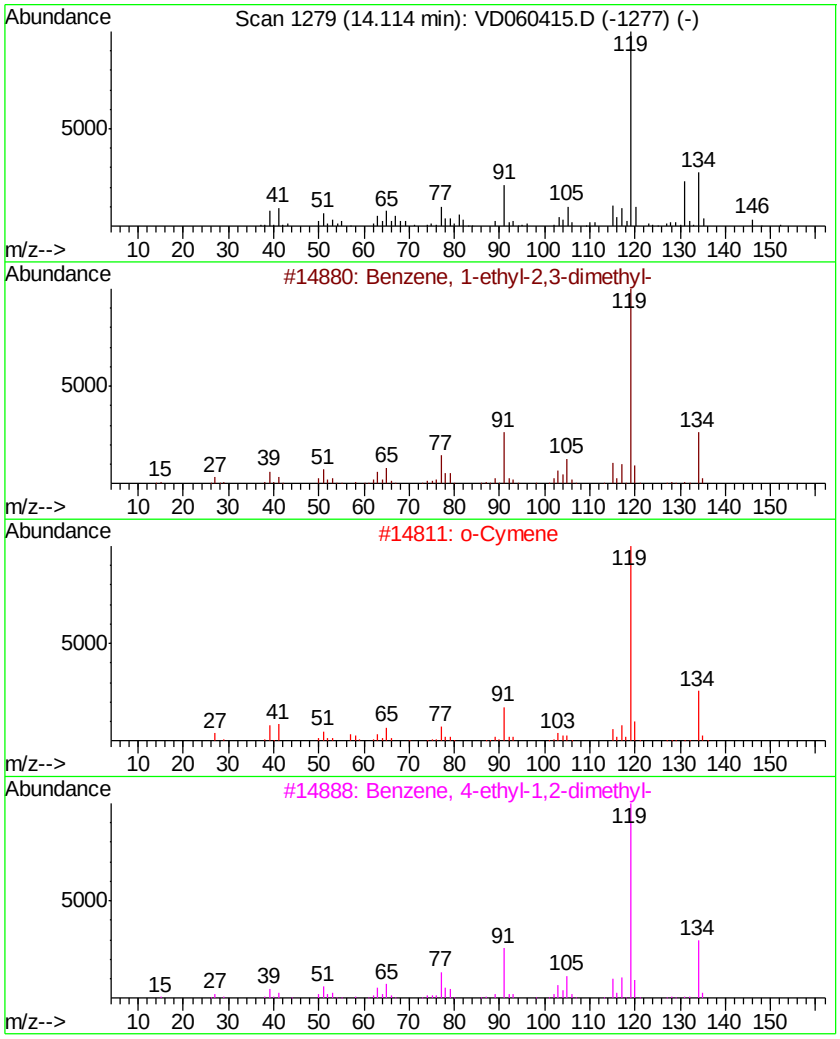
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

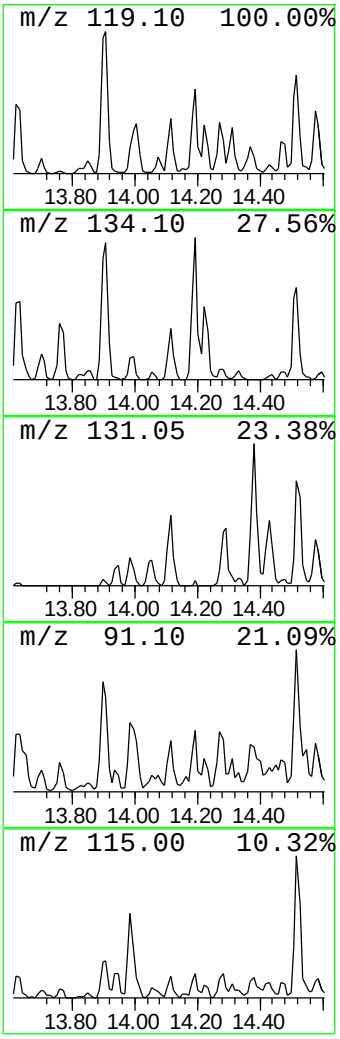
****

Peak Number 3 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.11	5.98 ug/l	436446	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	94
2	o-Cymene	134	C10H14	000527-84-4	94
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	94
4	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	94
5	p-Cymene	134	C10H14	000099-87-6	93





Library Search Compound Report

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ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

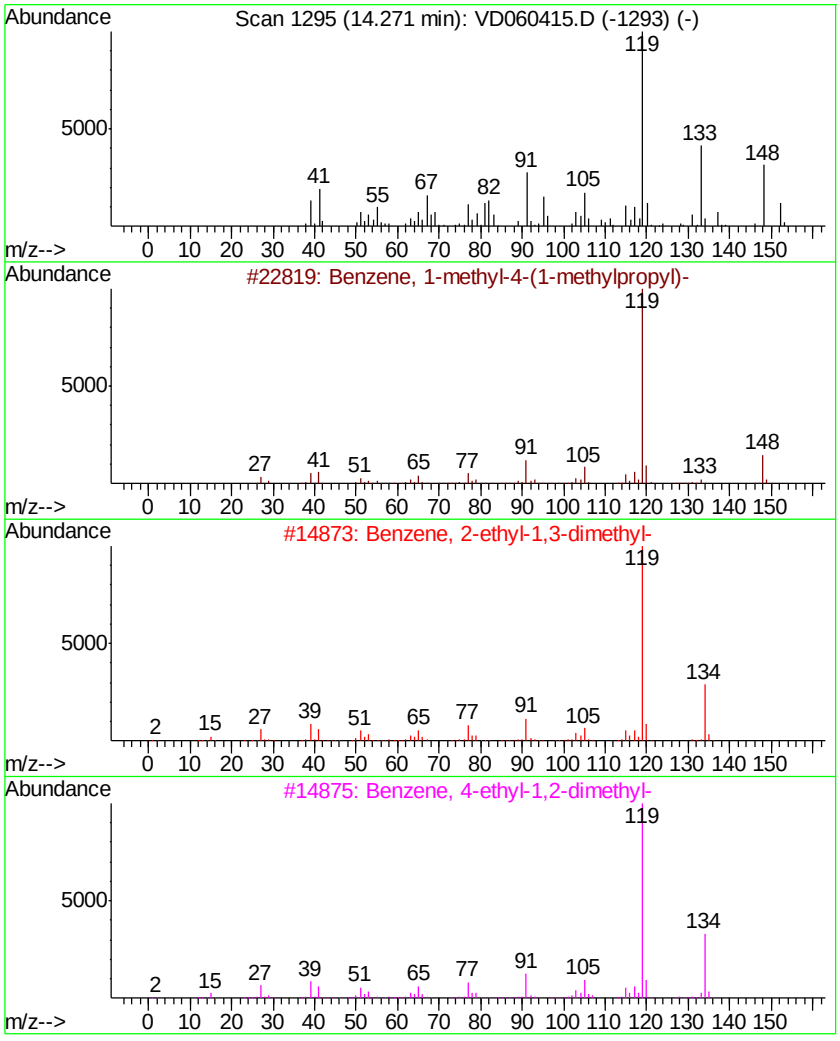
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

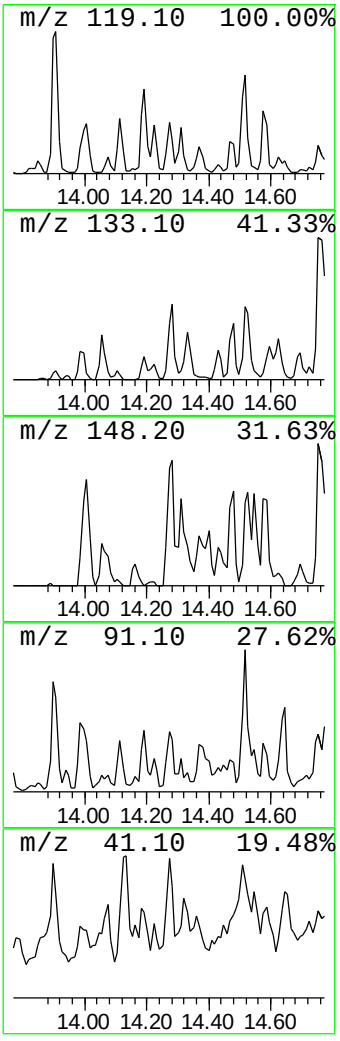
****

Peak Number 4 Benzene, 1-methyl-4-(1-meth... Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.27	6.19 ug/l	451974	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-4-(1-methylpro...	148	C11H16	001595-16-0	53
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	50
3	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	50
4	Benzene, 1-ethyl-2,3-dimethyl-	134	C10H14	000933-98-2	50
5	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	50





Library Search Compound Report

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Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

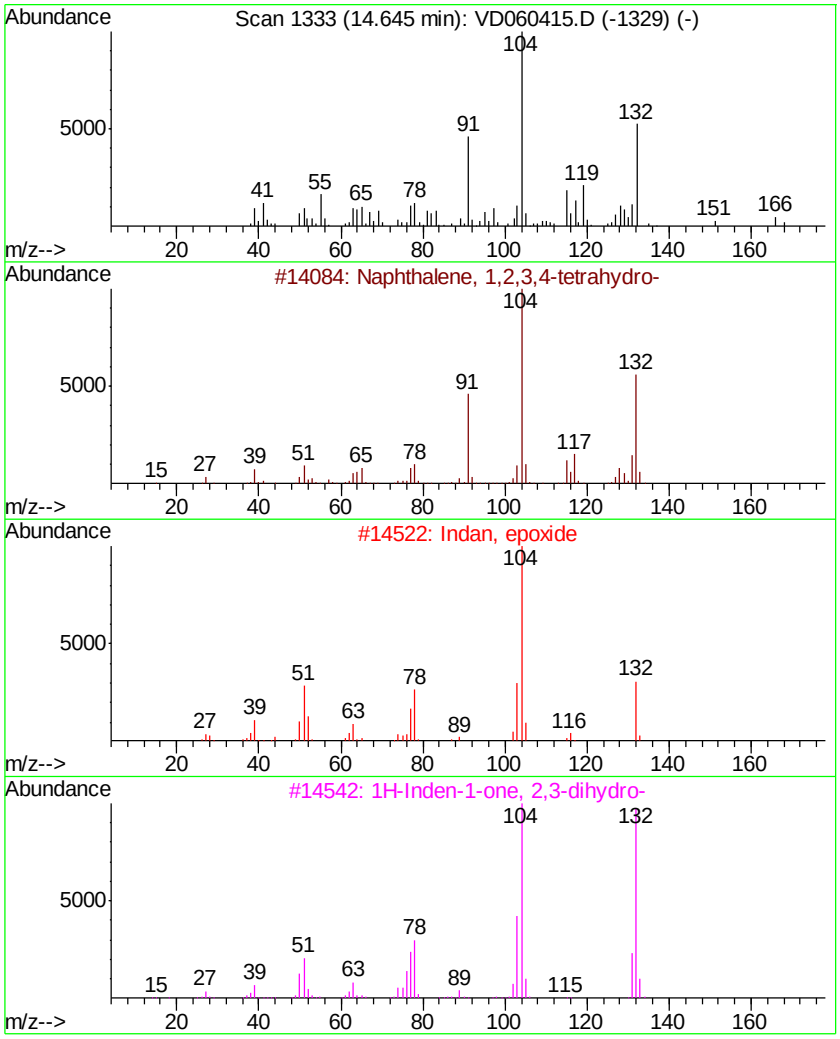
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

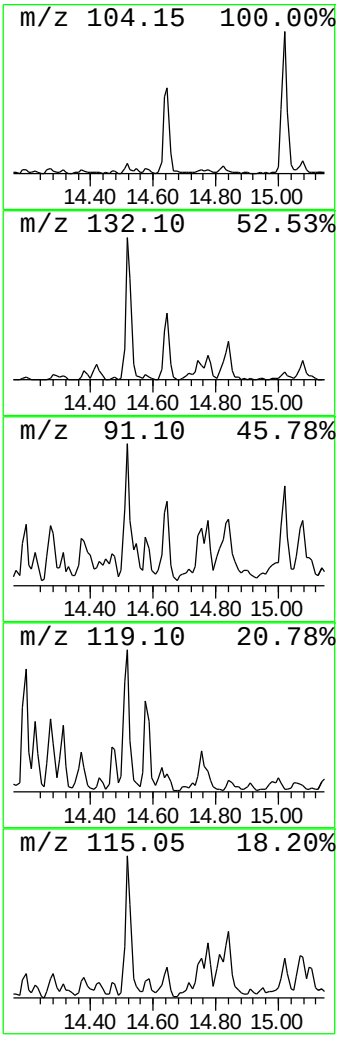
****

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

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.65	6.74 ug/l	492207	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	94
2	Indan, epoxide	132	C9H8O	000768-22-9	53
3	1H-Inden-1-one, 2,3-dihydro-	132	C9H8O	000083-33-0	38
4	2-Methylbenzyl cyanide	131	C9H9N	022364-68-7	35
5	Benzenebutanal	148	C10H12O	018328-11-5	35





Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
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Acq On : 19 Nov 2018 14:34
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Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

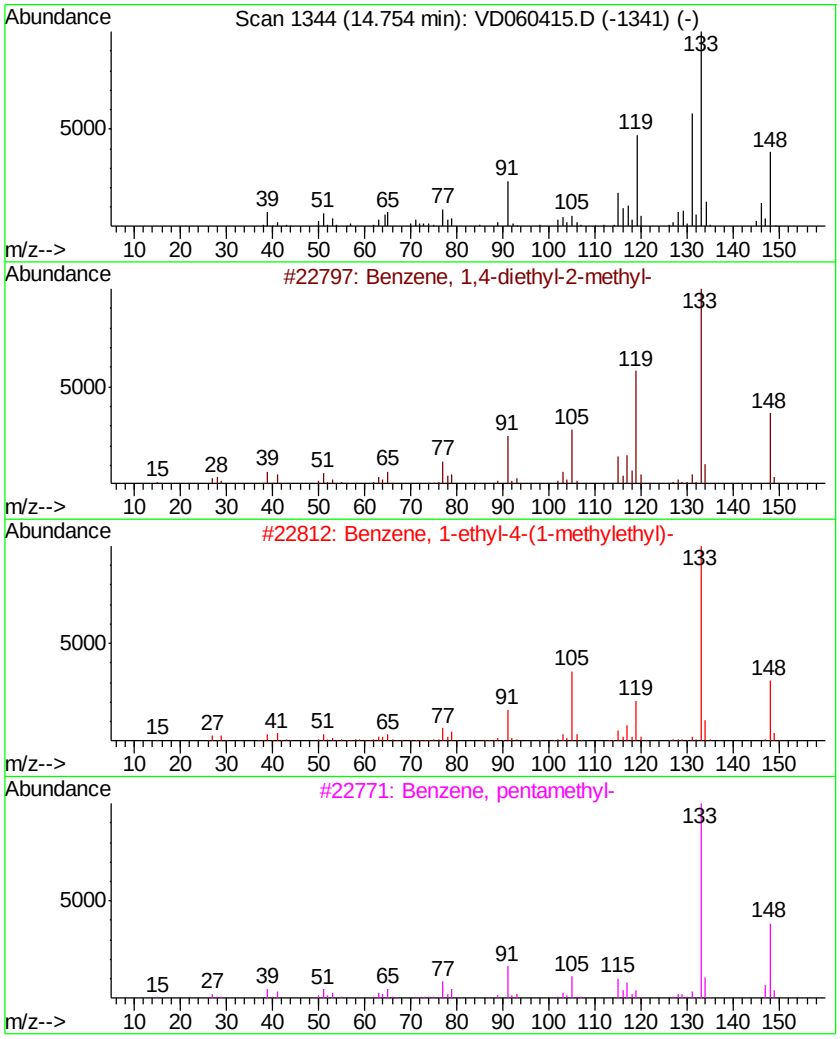
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

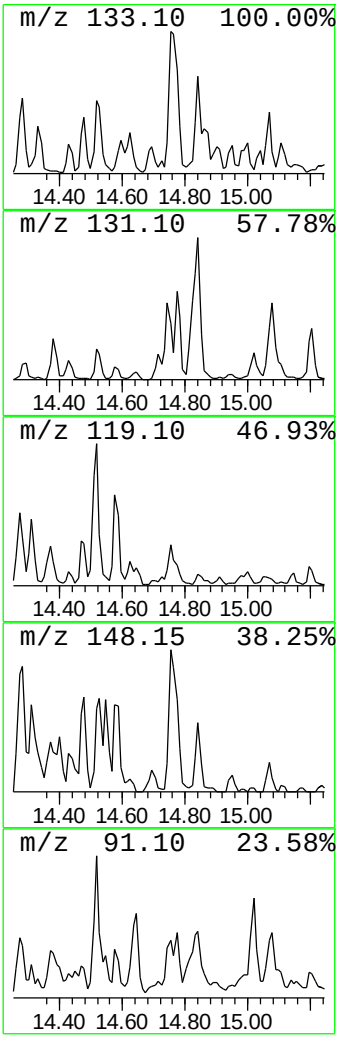
****

Peak Number 7 Benzene, 1,4-diethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.75	6.16 ug/l	449565	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1,4-diethyl-2-methyl-	148	C11H16	013632-94-5	68
2	Benzene, 1-ethyl-4-(1-methylethyl)-	148	C11H16	004218-48-8	52
3	Benzene, pentamethyl-	148	C11H16	000700-12-9	49
4	Benzene, 1,3-dimethyl-5-(1-methy...	148	C11H16	004706-90-5	49
5	3,4-Dimethylcumene	148	C11H16	1000370-34-1	47





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Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

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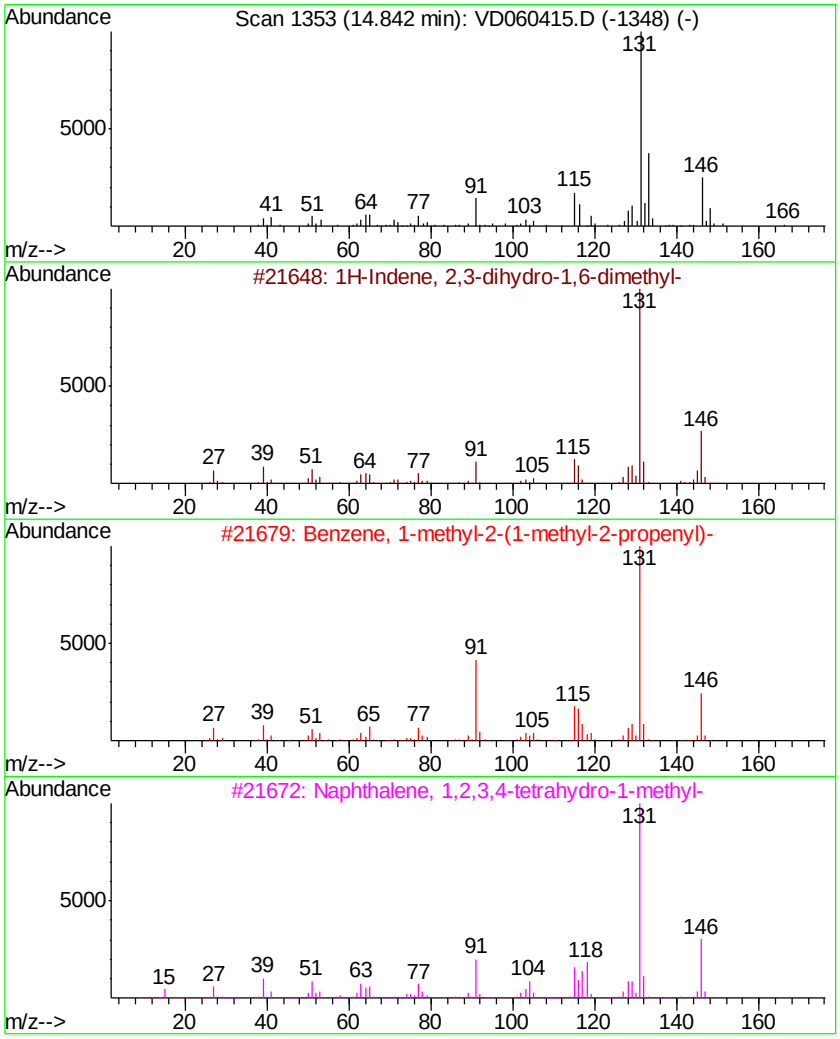
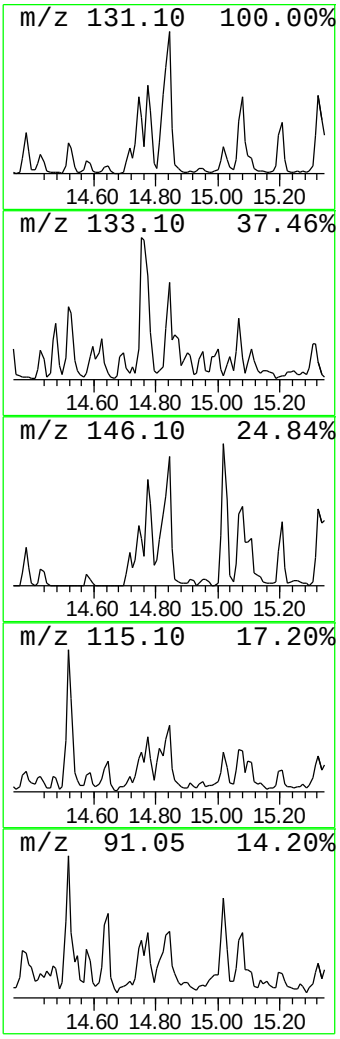
Peak Number 8 1H-Indene, 2,3-dihydro-1,6-... Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.84	13.75 ug/l	1003450	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1H-Indene, 2,3-dihydro-1,6-dimet...	146	C11H14	017059-48-2	93
2	Benzene, 1-methyl-2-(1-methyl-2-...	146	C11H14	097664-19-2	81
3	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	76
4	Benzene, (2-methyl-1-butenyl)-	146	C11H14	056253-64-6	76
5	Benzene, (1,1-dimethyl-2-propenyl)-	146	C11H14	018321-36-3	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

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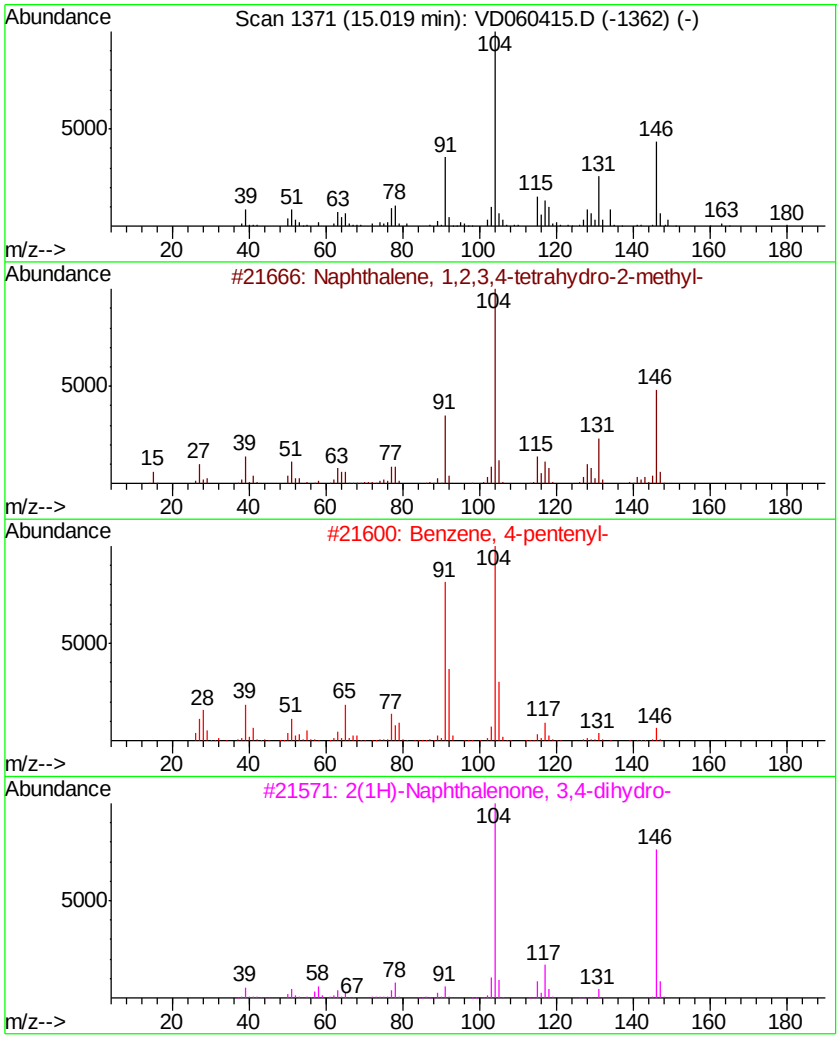
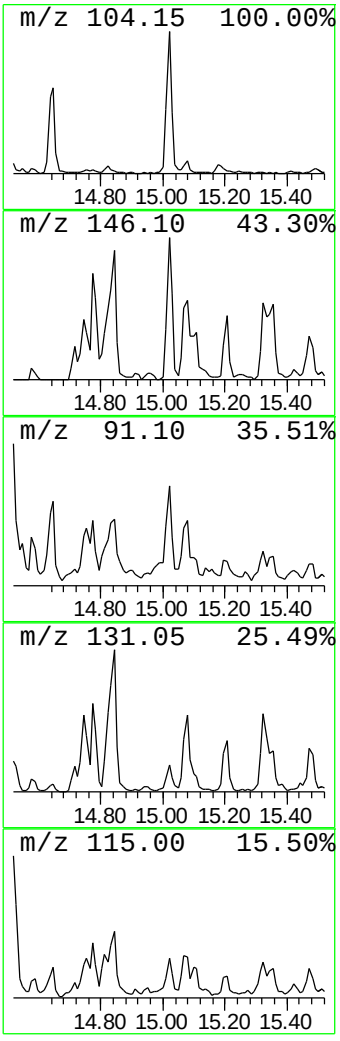
Peak Number 9 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.02	10.90 ug/l	795606	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	003877-19-8	94
2	Benzene, 4-pentenyl-	146	C11H14	001075-74-7	59
3	2(1H)-Naphthalenone, 3,4-dihydro-	146	C10H10O	000530-93-8	59
4	Naphth[1,2-b]oxirene, 1a,2,3,7b-...	146	C10H10O	002461-34-9	58
5	Naphthalene, 1,2,3,4-tetrahydro-	132	C10H12	000119-64-2	49





Data Path : Z:\VOASRV\HPCHEM1\MSVOA D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

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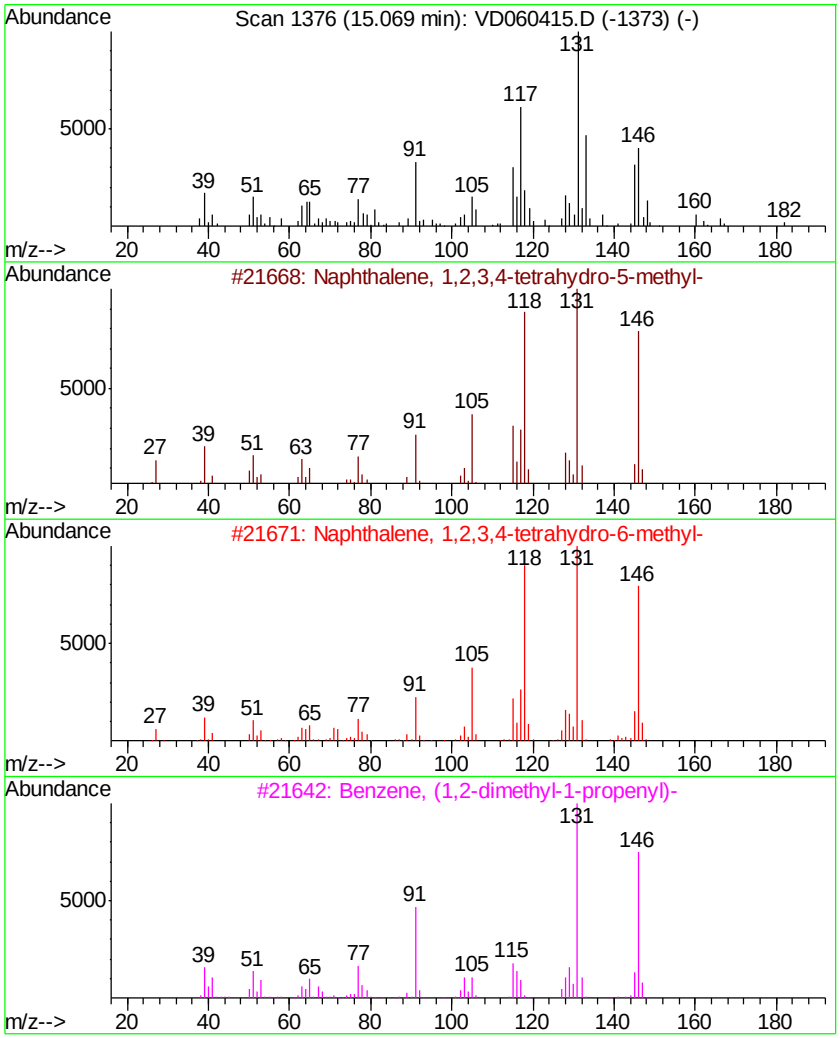
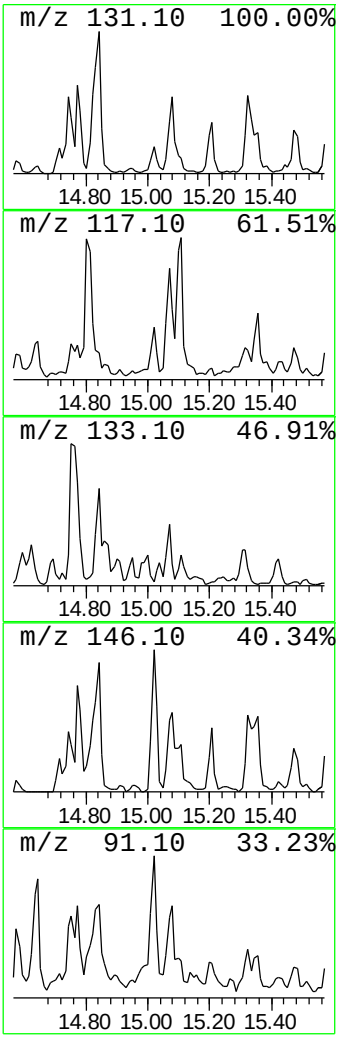
Peak Number 10 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.07	6.70 ug/l	489425	1,4-Dichlorobenzene-d4	13.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	83
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	64
3	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	60
4	1H-Indene, 2,3-dihydro-4,7-dimet...	146	C11H14	006682-71-9	60
5	Benzene, 1-methyl-4-(1-methyl-2-...	146	C11H14	097664-18-1	60





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_D\DATA\VD111918\
Data File : VD060415.D
Acq On : 19 Nov 2018 14:34
Operator : JC/SY
Sample : J5774-01
Misc : 7.74g/5ml/MSVOA_D/SOIL
ALS Vial : 6 Sample Multiplier: 1

Instrument : MSVOA_D
ClientSampleId : SU-04-111618

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_D\METHOD\82D110618S.M
Quant Title : SW846 8260

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

					Internal Standard			
TIC Top Hit name	RT	EstConc	Units	Response	#	RT	Resp	Conc
p-Cymene	13.90	7.3	ug/l	534363	4	13.38	3650070	50.0
1-Phenyl-1-butene	13.99	7.9	ug/l	578785	4	13.38	3650070	50.0
Benzene, 1-ethyl-...	14.11	6.0	ug/l	436446	4	13.38	3650070	50.0
Benzene, 1-methyl...	14.27	6.2	ug/l	451974	4	13.38	3650070	50.0
Naphthalene, 1,2,...	14.65	6.7	ug/l	492207	4	13.38	3650070	50.0
Benzene, 1,4-diet...	14.75	6.2	ug/l	449565	4	13.38	3650070	50.0
1H-Indene, 2,3-di...	14.84	13.8	ug/l	1003450	4	13.38	3650070	50.0
Naphthalene, 1,2,...	15.02	10.9	ug/l	795606	4	13.38	3650070	50.0
Naphthalene, 1,2,...	15.07	6.7	ug/l	489425	4	13.38	3650070	50.0