LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

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_N\METHODS\82N100418W.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	7.593	1790	1810	1821	rBV2	544966	1367905	13.34%	1.228%
2	7.667	1821	1833	1863	rVB	861184	2033645	19.83%	1.825%
3	8.030	1928	1946	1968	rBV	537403	1232290	12.02%	1.106%
4	8.278	1990	2023	2025	rBV4	41496	163213	1.59%	0.146%
5	8.587	2103	2119	2152	rBV	1139552	2364557	23.06%	2.122%
6	10.091	2571	2587	2616	rBV	2145645	3990241	38.91%	3.581%
7	10.718	2771	2782	2794	rVB	95151	162003	1.58%	0.145%
8	10.821	2798	2814	2827	rBV	60552	129948	1.27%	0.117%
9	11.008	2861	2872	2882	rBV2	207177	358978	3.50%	0.322%
10	11.124	2896	2908	2916	rBV3	158848	295693	2.88%	0.265%
11	11.214	2926	2936	2951	rVB3	200822	442832	4.32%	0.397%
12	11.300	2951	2963	2972	rBV2	197076	362228	3.53%	0.325%
13	11.410	2983	2997	3016	rBV	1616912	2952083	28.79%	2.650%
14	11.545	3030	3039	3045	rBV2	139963	226891	2.21%	0.204%
15	11.596	3045	3055	3064	rVV5	226952	459698	4.48%	0.413%
16	11.657	3064	3074	3081	rBV	600722	1075466	10.49%	0.965%
17	11.699	3082	3087	3097	rVB2	428530	657156	6.41%	0.590%
18	11.760	3097	3106	3112	rBV3	152568	276534	2.70%	0.248%
19	11.808	3113	3121	3126	rBV5	62667	118754	1.16%	0.107%
20	11.853	3127	3135	3143	rVB4	383590	594443	5.80%	0.534%
21	11.927	3143	3158	3169	rBV6	1216373	2876218	28.05%	2.581%
22	11.979	3170	3174	3179	rBV3	193208	208762	2.04%	0.187%
23	12.046	3186	3195	3204	rVB	2366918	3374820	32.91%	3.029%
24	12.120	3209	3218	3223	rBV3	1137750	1972998	19.24%	1.771%
25	12.159	3224	3230	3241	rVB5	2001732	3680911	35.89%	3.304%
26	12.297	3264	3273	3279	rBV6	955286	1690128	16.48%	1.517%
27	12.336	3280	3285	3294	rVB	2059770	3012302	29.37%	2.704%
28	12.403	3294	3306	3317	rVB3	2092332	3905653	38.09%	3.505%
29	12.493	3318	3334	3340	rBV5	737184	1942398	18.94%	1.743%
30	12.564	3348	3356	3371	rVB4	2234718	4652341	45.37%	4.176%
31	12.644	3371	3381	3389	rBV7	1069457	2339686	22.82%	2.100%
32	12.731	3390	3408	3417	rBV5	3028292	8097192	78.96%	7.267%
33	12.869	3444	3451	3458	rBV5	1173424	1752722	17.09%	1.573%
34	12.963	3473	3480	3493	rVB	6120092	9701509	94.60%	8.707%

LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

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_N\METHODS\82N100418W.M
Title : SW846 8260

35	13.040	3494	3504	3513	rBV	6444833	10255006	100.00%	9.204%
36	13.091	3514	3520	3529	rVB3	1007137	1489364	14.52%	1.337%
37	13.146	3530	3537	3541	rBV3	1276839	1894743	18.48%	1.701%
38	13.239	3552	3566	3574	rBV3	2737323	5526169	53.89%	4.960%
39	13.284	3575	3580	3588	rVB4	1663595	2318833	22.61%	2.081%
40	13.590	3667	3675	3688	rBV4	1281509	2638127	25.73%	2.368%
41	13.664	3689	3698	3708	rBV5	2246337	4160082	40.57%	3.734%
42	13.834	3746	3751	3760	rVB4	1599299	2272350	22.16%	2.039%
43	13.889	3761	3768	3777	rVB	1892080	2923267	28.51%	2.624%
44	13.995	3792	3801	3810	rVB4	1239052	2097356	20.45%	1.882%
45	14.127	3838	3842	3848	rBV2	294608	316262	3.08%	0.284%
46	14.178	3850	3858	3866	rBV5	721912	1356452	13.23%	1.217%
47	14.249	3874	3880	3888	rVB	885259	1235374	12.05%	1.109%
48	14.294	3888	3894	3900	rBV3	441480	604362	5.89%	0.542%
49	14.332	3901	3906	3913	rVB3	496322	642130	6.26%	0.576%
50	14.525	3960	3966	3974	rVB2	300814	420988	4.11%	0.378%
51	14.583	3974	3984	3996	rVV2	976707	1881091	18.34%	1.688%
52	14.644	3997	4003	4015	rVB2	250037	421901	4.11%	0.379%
53	14.741	4027	4033	4042	rVB	248238	358408	3.49%	0.322%
54	14.853	4062	4068	4073	rBV5	107017	135453	1.32%	0.122%

Sum of corrected areas: 111419916

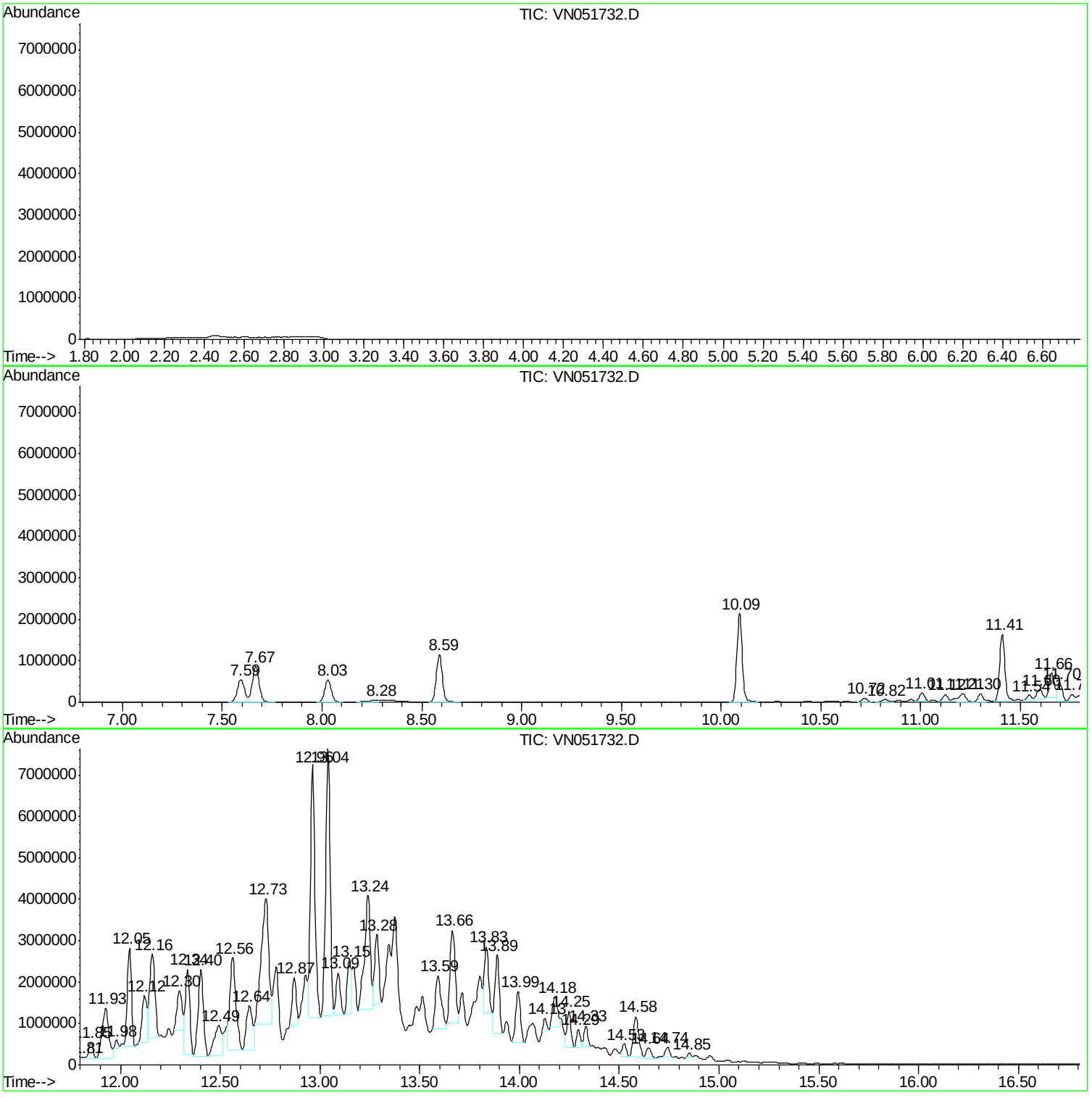
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument :
MSVOA_N
ClientSampleId :
EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

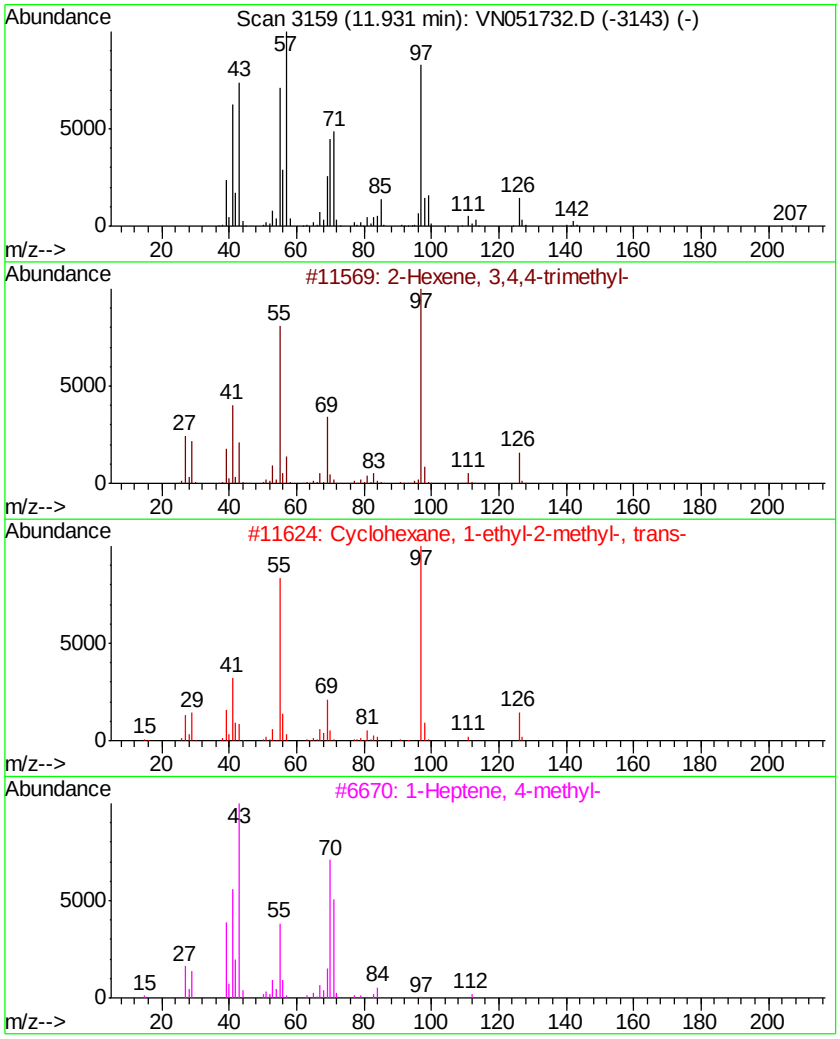
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

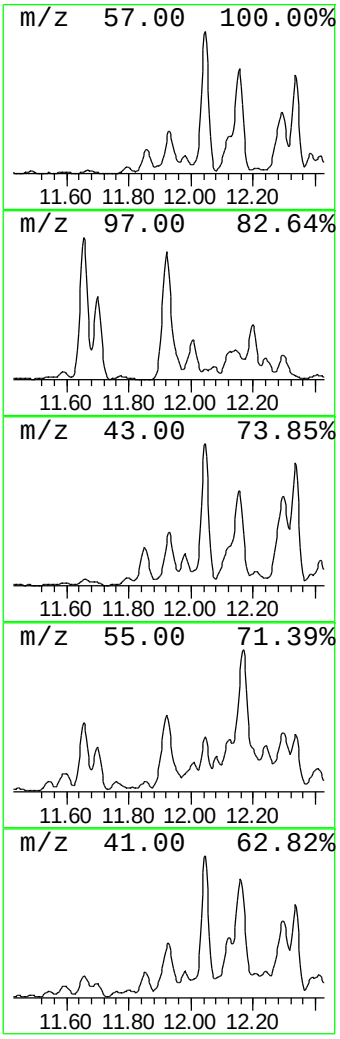
****

Peak Number 1 2-Hexene, 3,4,4-trimethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.93	48.72 ug/l	2876220	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	2-Hexene, 3,4,4-trimethyl-	126	C9H18	053941-19-8	43
2	Cyclohexane, 1-ethyl-2-methyl-, ...	126	C9H18	004923-78-8	43
3	1-Heptene, 4-methyl-	112	C8H16	013151-05-8	38
4	4-Octen-3-one	126	C8H14O	014129-48-7	38
5	3,5-Dimethyl-3-heptene	126	C9H18	059643-68-4	35





Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

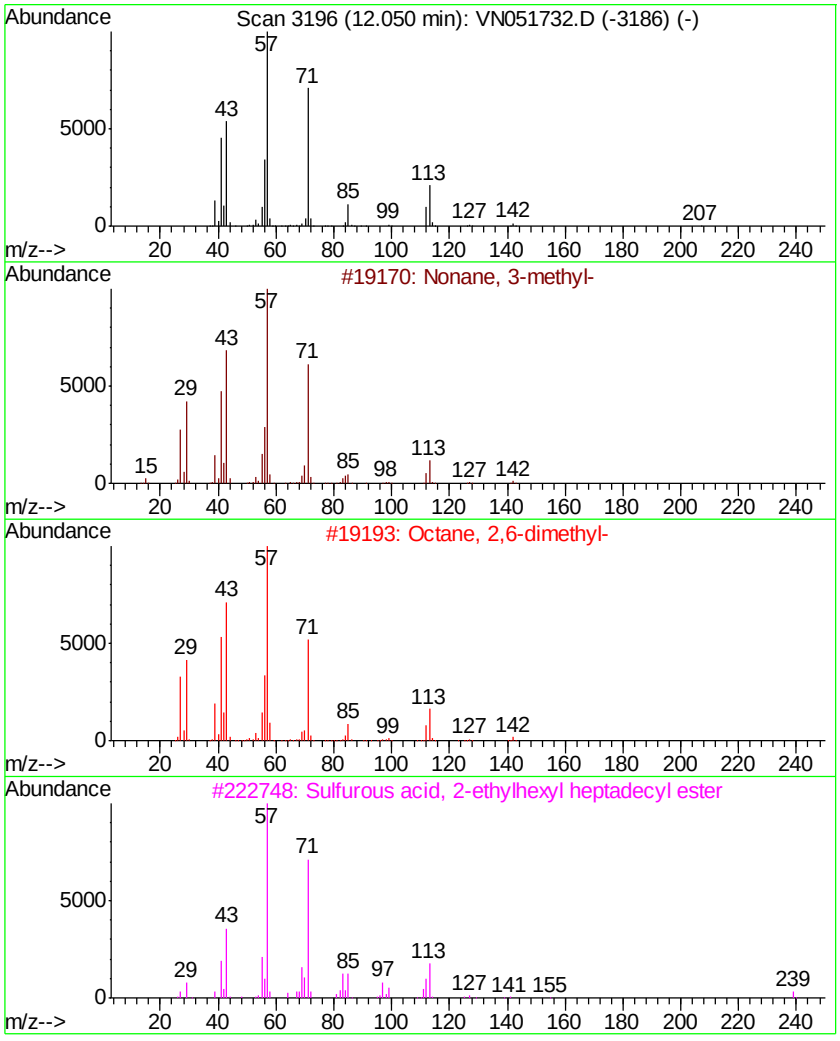
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

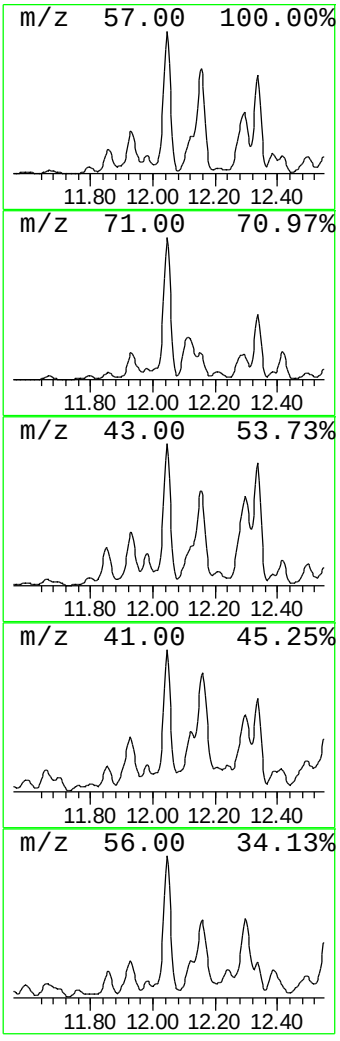
****

Peak Number 2 Nonane, 3-methyl- Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.05	57.16 ug/l	3374820	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 3-methyl-	142	C10H22	005911-04-6	91
2	Octane, 2,6-dimethyl-	142	C10H22	002051-30-1	90
3	Sulfurous acid, 2-ethylhexyl hep...	432	C25H52O3S	1000309-20-0	72
4	Octane, 3,6-dimethyl-	142	C10H22	015869-94-0	72
5	Octane, 1,1'-oxybis-	242	C16H34O	000629-82-3	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

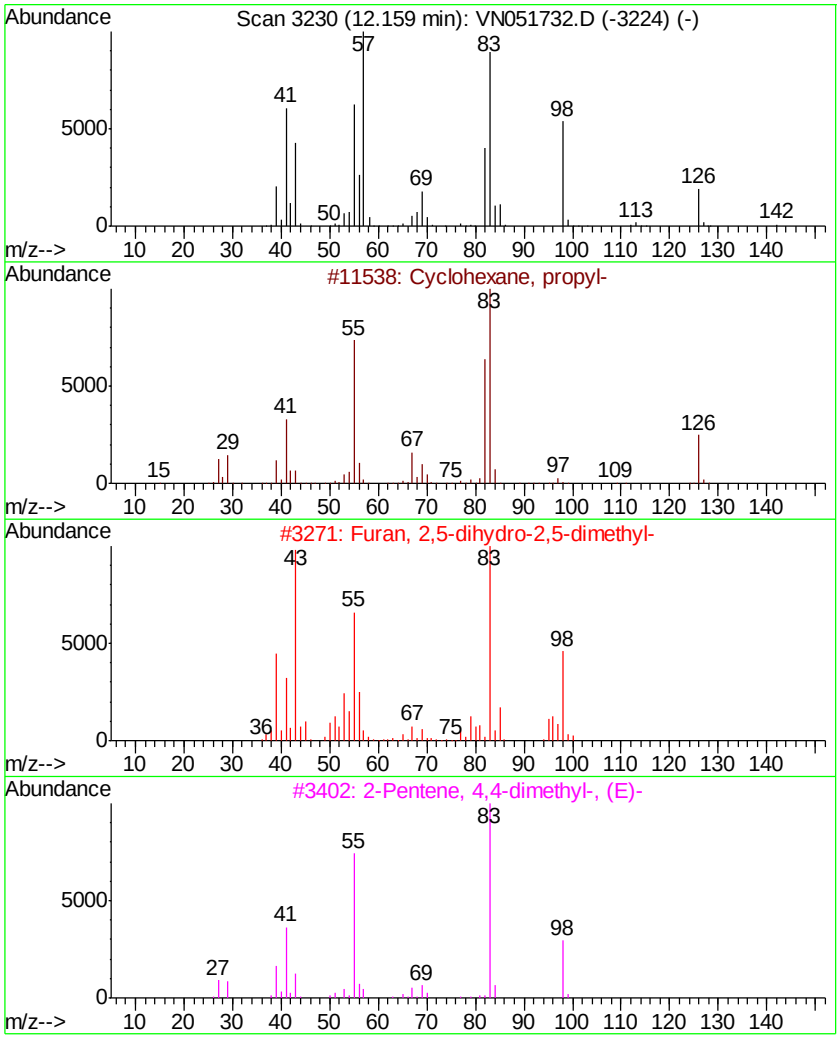
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

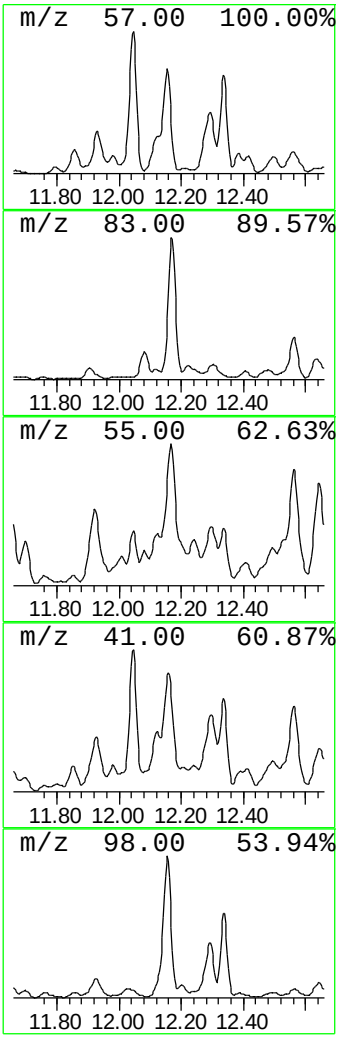
****

Peak Number 3 Cyclohexane, propyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.16	62.34 ug/l	3680910	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, propyl-	126	C9H18	001678-92-8	49
2	Furan, 2,5-dihydro-2,5-dimethyl-	98	C6H10O	059242-27-2	47
3	2-Pentene, 4,4-dimethyl-, (E)-	98	C7H14	000690-08-4	47
4	2-Pentene, 3,4-dimethyl-, (E)-	98	C7H14	004914-92-5	47
5	Cyclohexane, (1-methylethyl)-	126	C9H18	000696-29-7	43





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

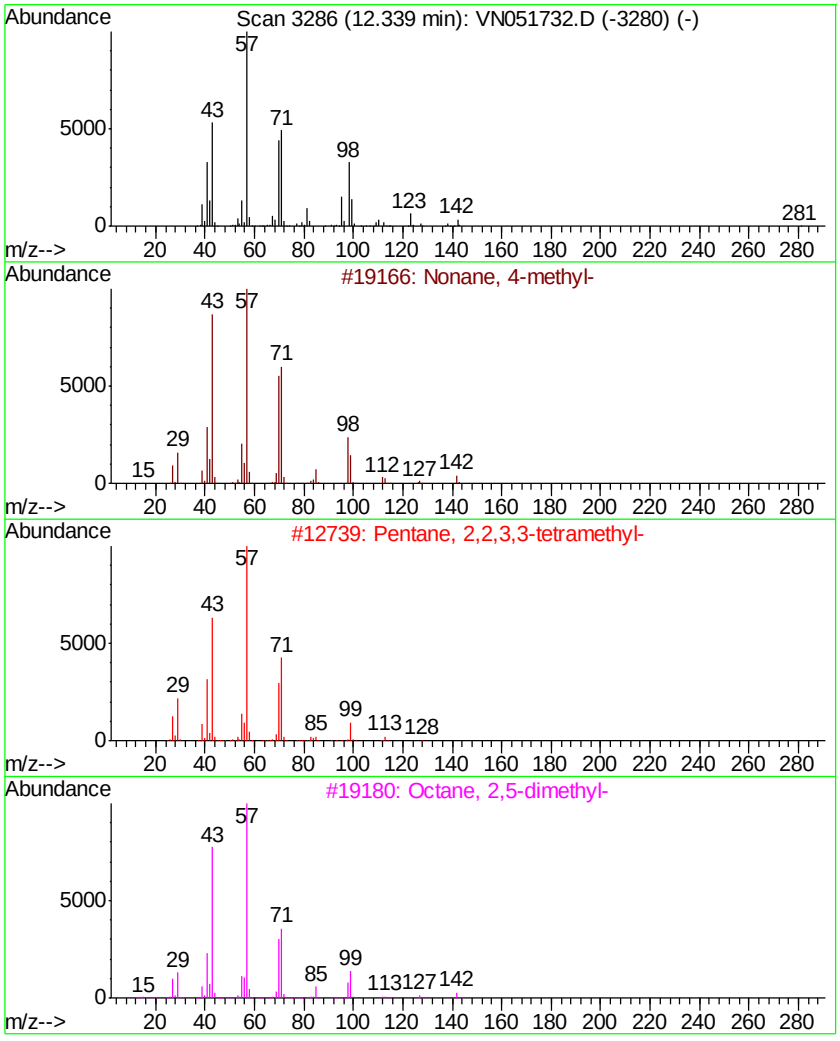
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

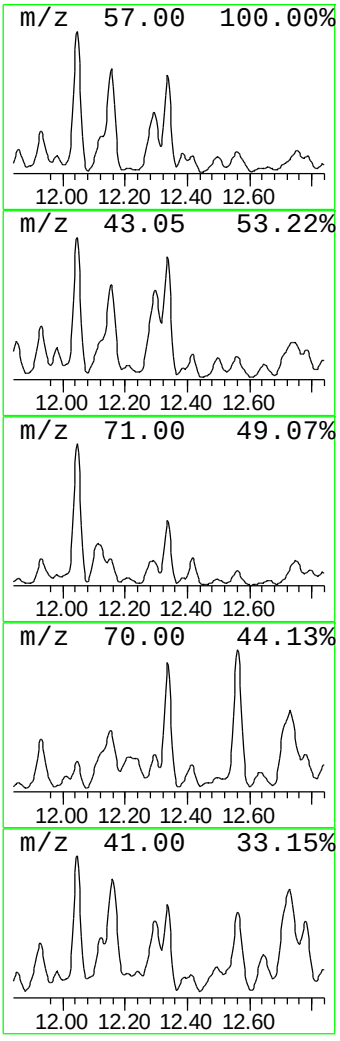
****

Peak Number 4 Nonane, 4-methyl- Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.34	51.02 ug/l	3012300	Chlorobenzene-d5	11.41

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 4-methyl-	142	C10H22	017301-94-9	64
2	Pentane, 2,2,3,3-tetramethyl-	128	C9H20	007154-79-2	53
3	Octane, 2,5-dimethyl-	142	C10H22	015869-89-3	50
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	47
5	Heptane, 2,3,4-trimethyl-	142	C10H22	052896-95-4	45





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

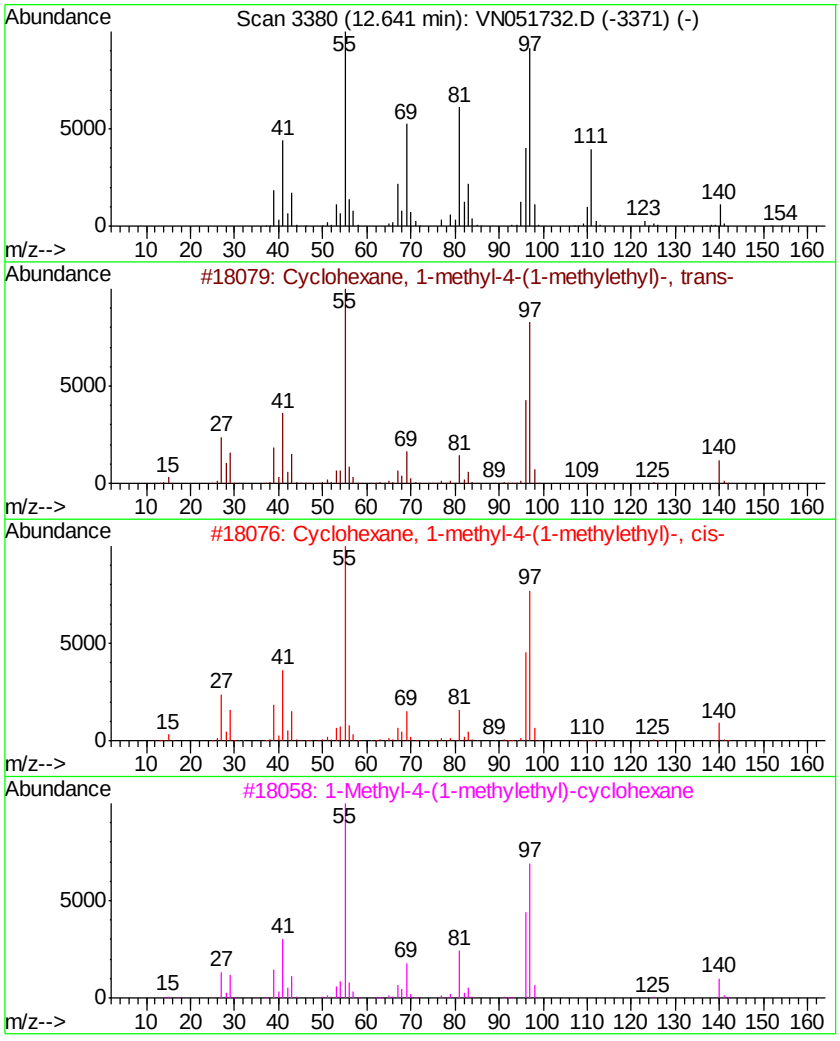
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

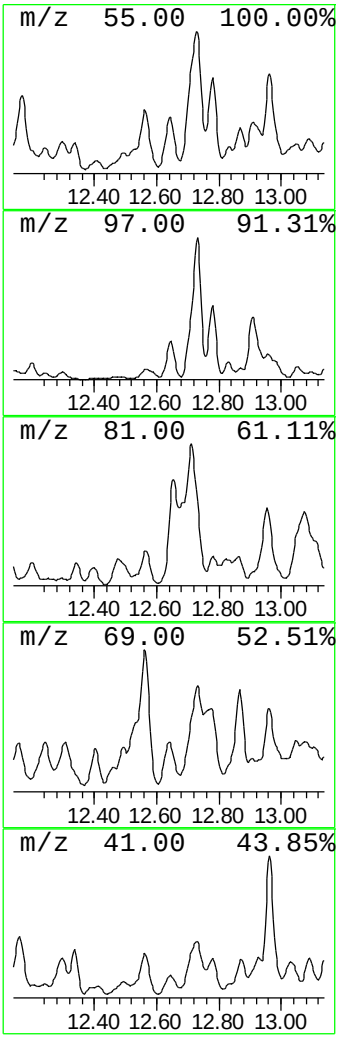
****

Peak Number 5 Cyclohexane, 1-methyl-4-(1-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.64	50.45 ug/l	2339690	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	001678-82-6	46
2	Cyclohexane, 1-methyl-4-(1-methy...	140	C10H20	006069-98-3	46
3	1-Methyl-4-(1-methylethyl)-cyclo...	140	C10H20	000099-82-1	45
4	Cyclohexene, 1-ethyl-	110	C8H14	001453-24-3	42
5	Cyclopentane, 1-hydroxymethyl-1,...	128	C8H16O	1000156-73-8	38





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

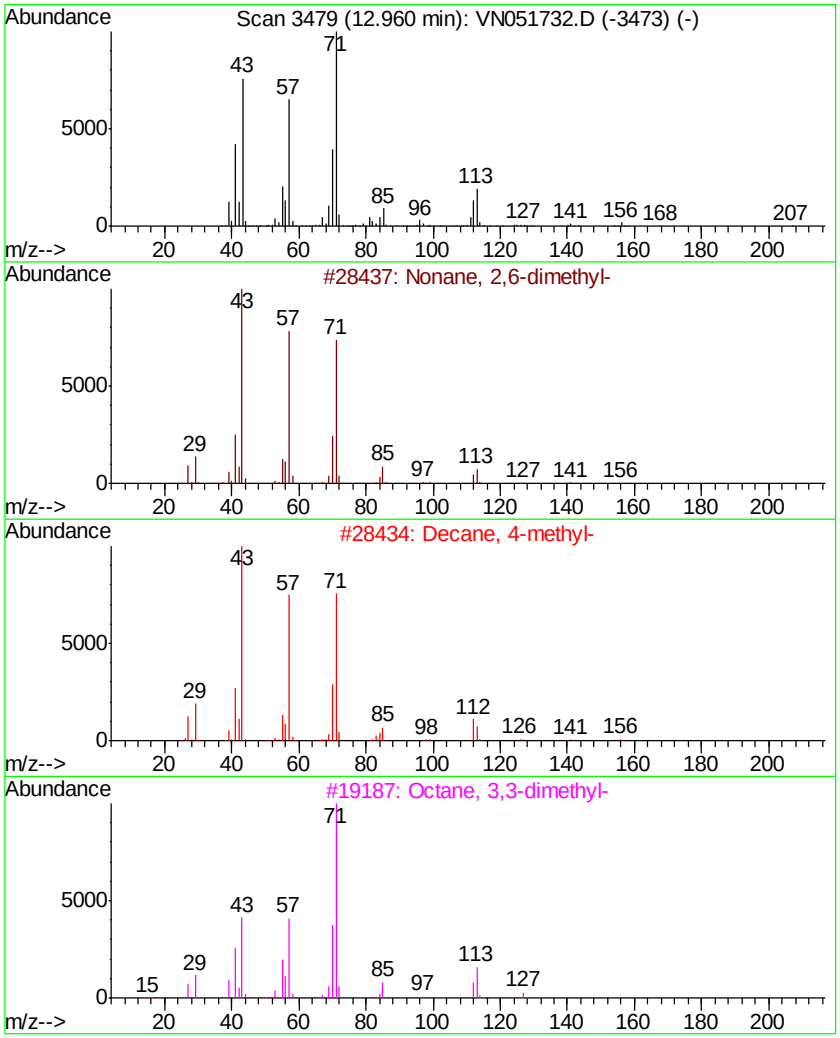
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

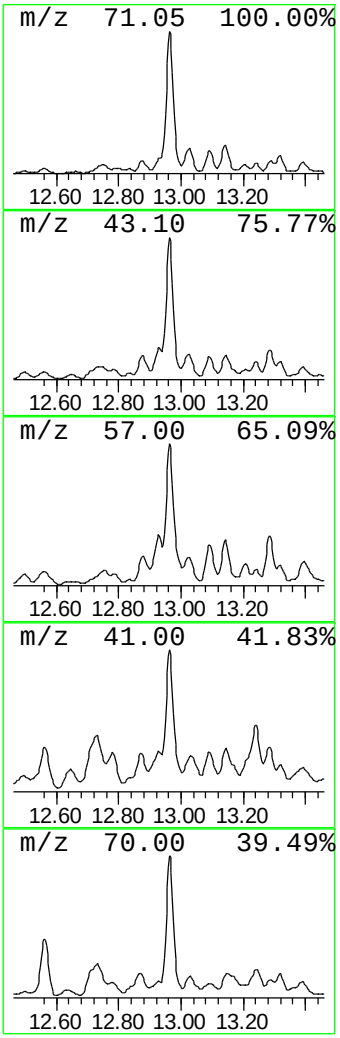
****

Peak Number 6 Nonane, 2,6-dimethyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.96	209.19 ug/l	9701510	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Nonane, 2,6-dimethyl-	156	C11H24	017302-28-2	87
2	Decane, 4-methyl-	156	C11H24	002847-72-5	83
3	Octane, 3,3-dimethyl-	142	C10H22	004110-44-5	83
4	Heptane, 3,3,5-trimethyl-	142	C10H22	007154-80-5	78
5	Sulfurous acid, 2-ethylhexyl iso...	250	C12H26O3S	1000309-18-7	59





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

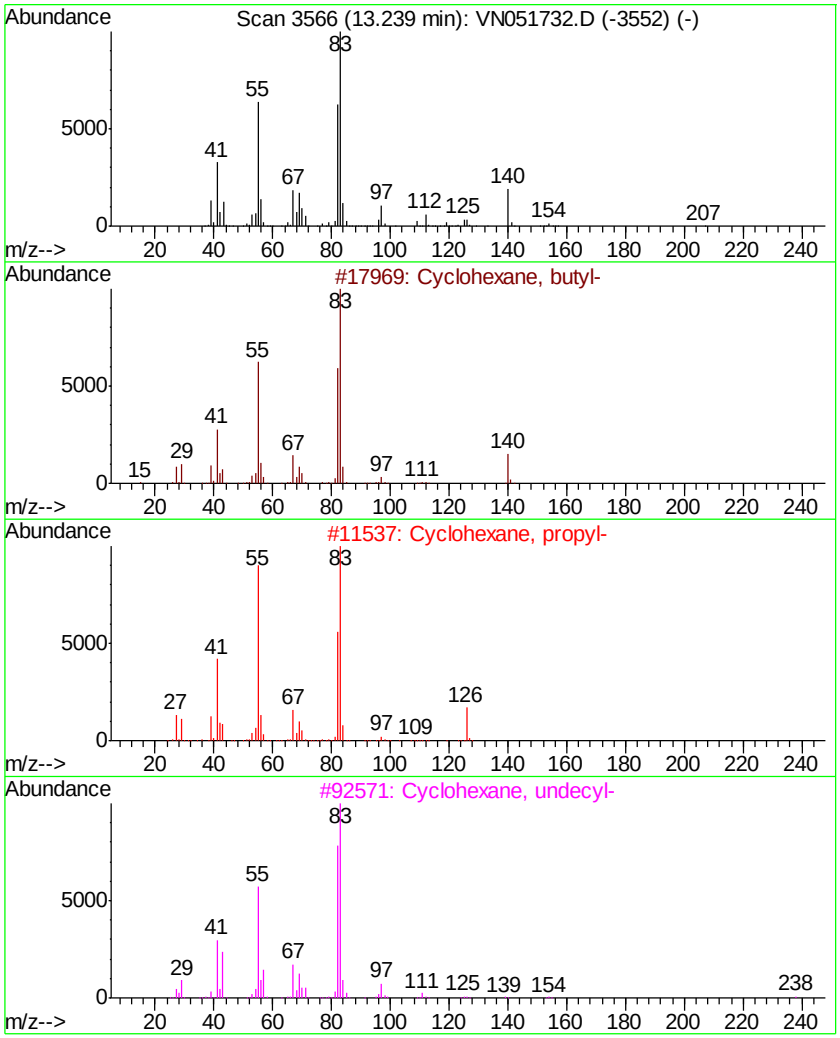
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

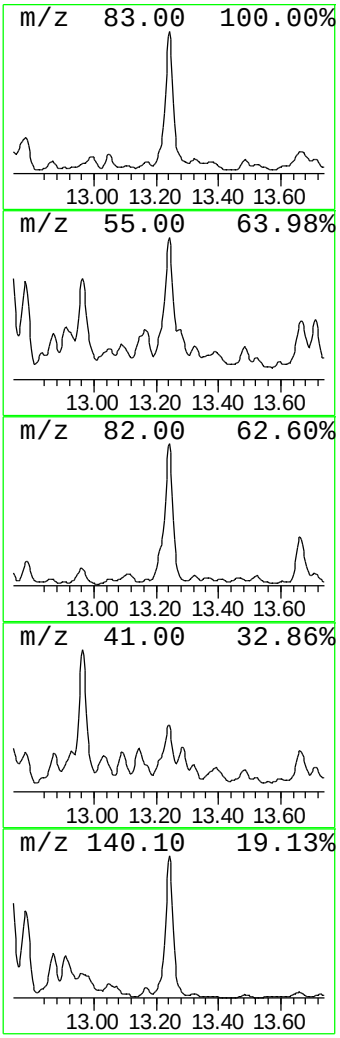
****

Peak Number 7 Cyclohexane, butyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.24	119.16 ug/l	5526170	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclohexane, butyl-	140	C10H20	001678-93-9	93
2	Cyclohexane, propyl-	126	C9H18	001678-92-8	81
3	Cyclohexane, undecyl-	238	C17H34	054105-66-7	80
4	Cyclohexane, pentyl-	154	C11H22	004292-92-6	74
5	Cyclohexane, ethyl-	112	C8H16	001678-91-7	72





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

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

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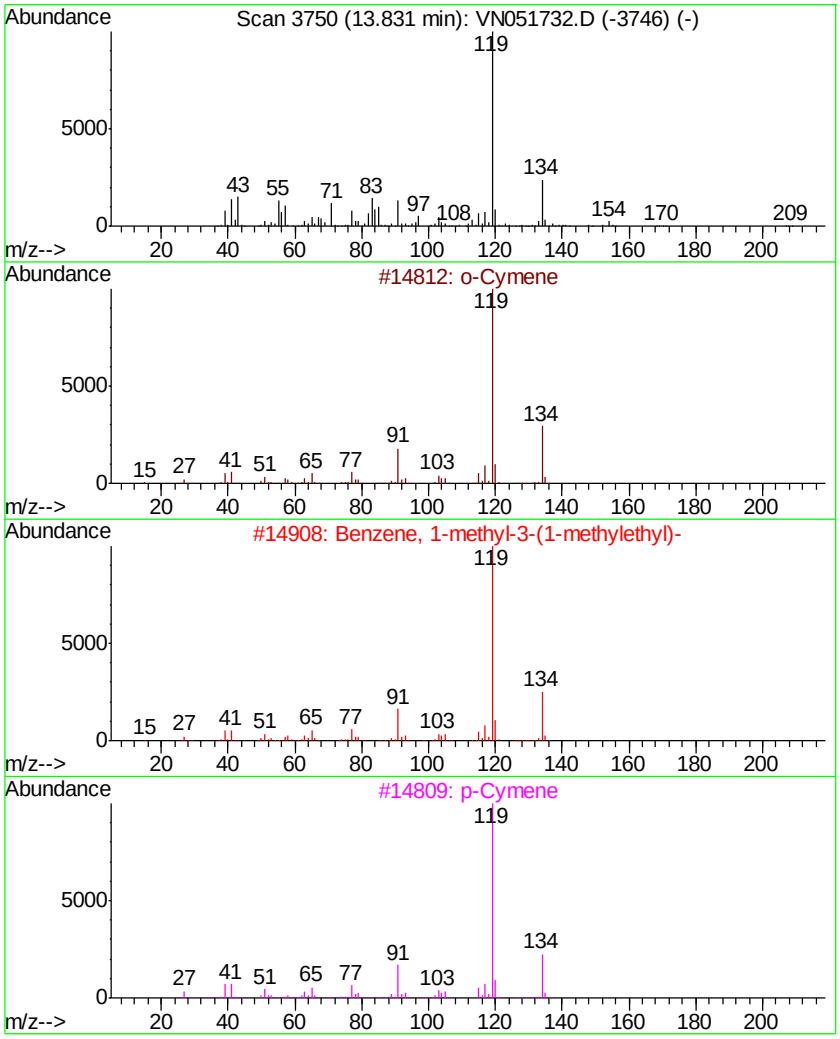
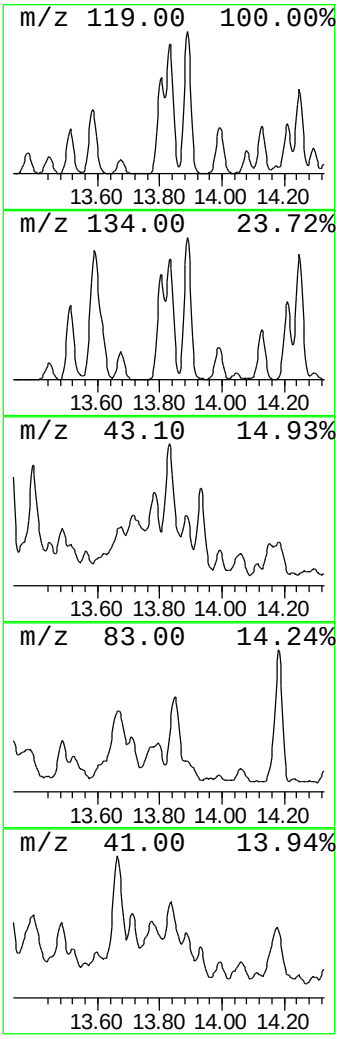
Peak Number 8 o-Cymene Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.83	49.00 ug/l	2272350	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	o-Cymene	134	C10H14	000527-84-4	92
2	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	92
3	p-Cymene	134	C10H14	000099-87-6	91
4	Benzene, 1,2,4,5-tetramethyl-	134	C10H14	000095-93-2	83
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	83





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

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

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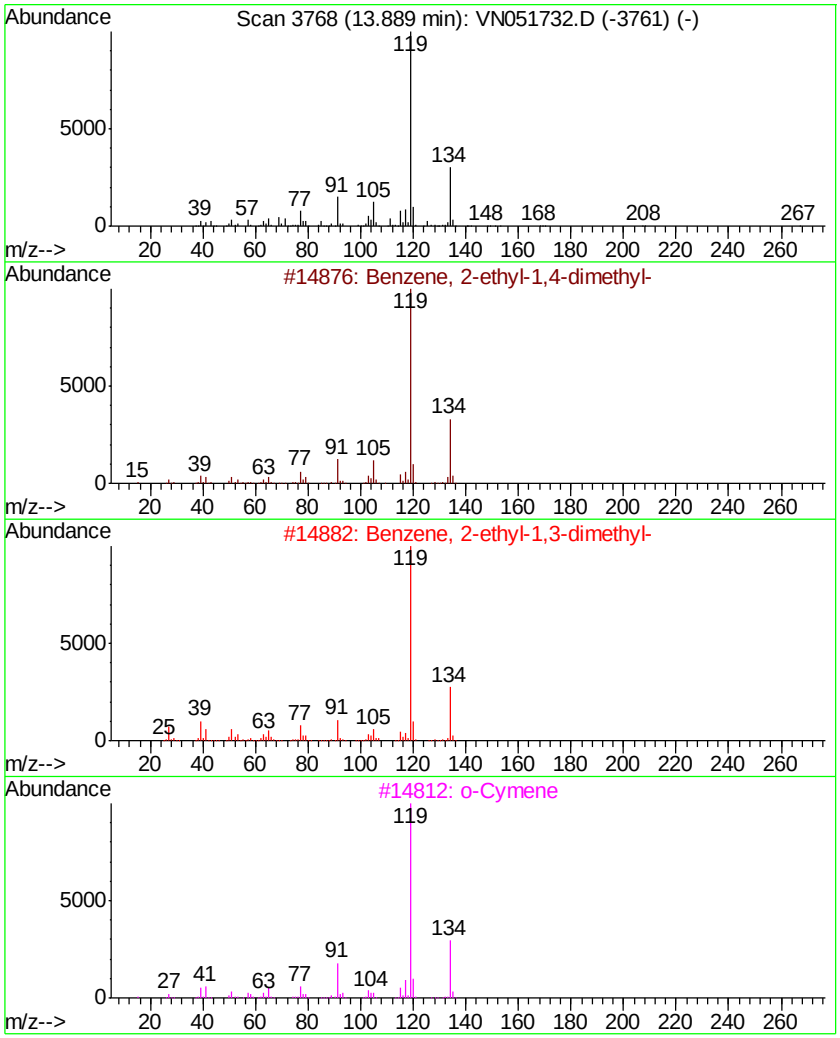
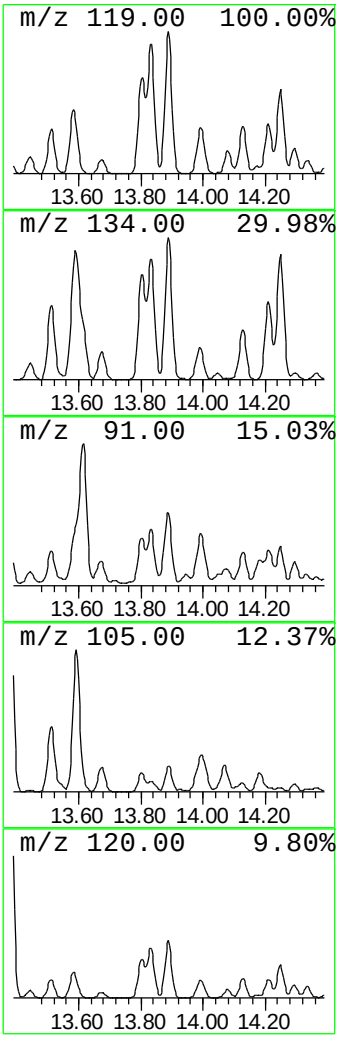
Peak Number 9 Benzene, 2-ethyl-1,4-dimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	63.03 ug/l	2923270	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	95
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	o-Cymene	134	C10H14	000527-84-4	95
4	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	95
5	Benzene, 1-methyl-3-(1-methyleth...	134	C10H14	000535-77-3	94





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.M
Quant Title : SW846 8260

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

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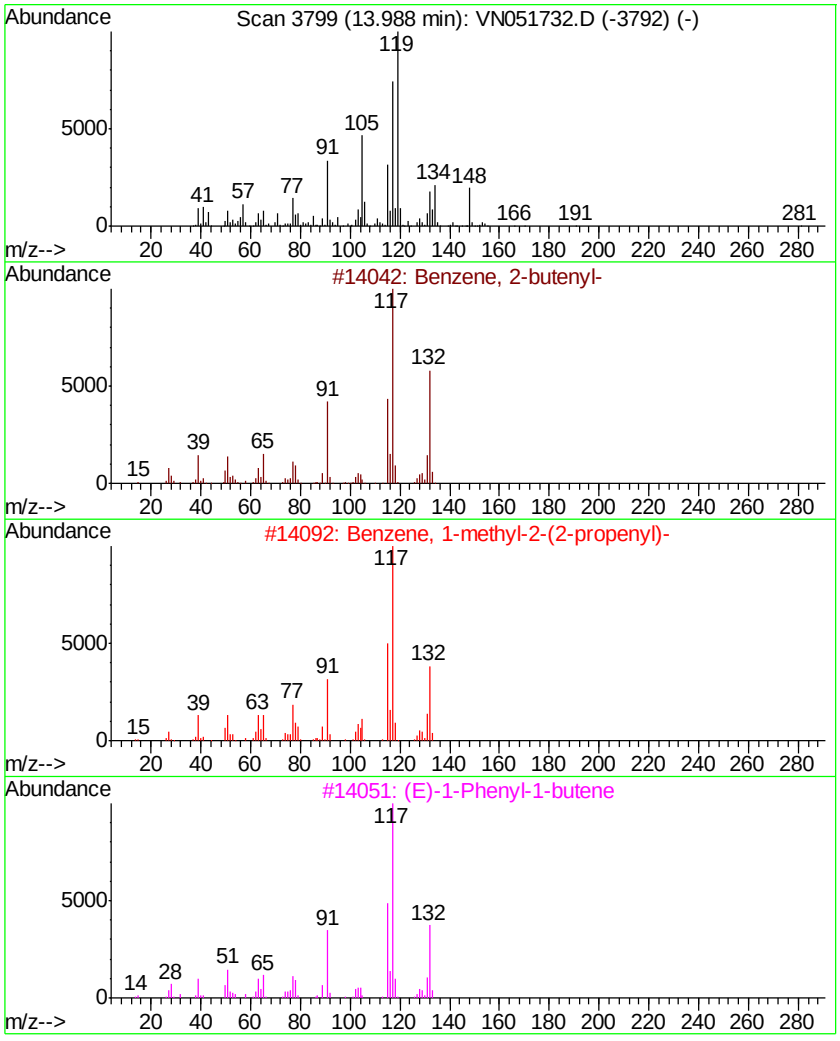
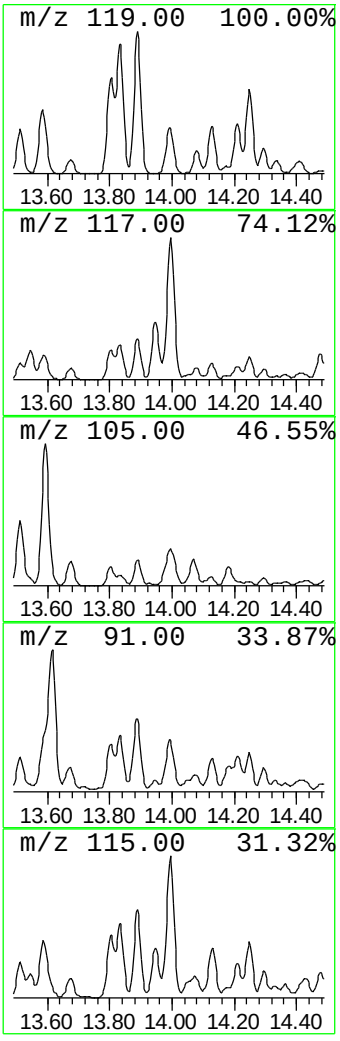
Peak Number 10 Benzene, 2-butenyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	45.22 ug/l	2097360	1,4-Dichlorobenzene-d4	13.35

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-butenyl-	132	C10H12	001560-06-1	86
2	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	70
3	(E)-1-Phenyl-1-butene	132	C10H12	001005-64-7	64
4	Benzene, (1-methyl-1-propenyl)-,...	132	C10H12	000768-00-3	50
5	Benzene, (2-methyl-2-propenyl)-	132	C10H12	003290-53-7	50





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN100818\
Data File : VN051732.D
Acq On : 8 Oct 2018 13:45
Operator : MD\SY
Sample : J5165-06 10X
Misc : 5.89g/5mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 15 Sample Multiplier: 28

Instrument : MSVOA_N
ClientSampleId : EP1-MW-H-092718(7-10)

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N100418W.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
2-Hexene, 3,4,4-t...	11.93	48.7	ug/l	2876220	3	11.41	2952080	50.0
Nonane, 3-methyl-	12.05	57.2	ug/l	3374820	3	11.41	2952080	50.0
Cyclohexane, propyl-	12.16	62.3	ug/l	3680910	3	11.41	2952080	50.0
Nonane, 4-methyl-	12.34	51.0	ug/l	3012300	3	11.41	2952080	50.0
Cyclohexane, 1-me...	12.64	50.5	ug/l	2339690	4	13.35	2318830	50.0
Nonane, 2,6-dimet...	12.96	209.2	ug/l	9701510	4	13.35	2318830	50.0
Cyclohexane, butyl-	13.24	119.2	ug/l	5526170	4	13.35	2318830	50.0
o-Cymene	13.83	49.0	ug/l	2272350	4	13.35	2318830	50.0
Benzene, 2-ethyl-...	13.89	63.0	ug/l	2923270	4	13.35	2318830	50.0
Benzene, 2-butenyl-	13.99	45.2	ug/l	2097360	4	13.35	2318830	50.0