LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

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\82N070820W.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	4.545	852	862	894	rVB2	135833	452400	2.31%	0.363%
2	5.005	977	1005	1035	rBV2	166936	599884	3.06%	0.481%
3	6.474	1439	1462	1478	rBV3	83727	268197	1.37%	0.215%
4	6.580	1478	1495	1514	rVB3	161969	488238	2.49%	0.392%
5	7.349	1721	1734	1753	rVB4	69619	199701	1.02%	0.160%
6	7.551	1775	1797	1802	rBV	145567	346925	1.77%	0.278%
7	7.622	1806	1819	1832	rVV5	384279	1227856	6.27%	0.985%
8	7.706	1832	1845	1868	rVB	352910	969916	4.95%	0.778%
9	7.837	1868	1886	1901	rBV	304891	751618	3.84%	0.603%
10	8.001	1918	1937	1958	rBV3	890712	2248620	11.48%	1.805%
11	8.124	1959	1975	1982	rBV4	307419	782424	3.99%	0.628%
12	8.265	2011	2019	2039	rVB2	112941	274044	1.40%	0.220%
13	8.548	2089	2107	2129	rBV3	616151	1491906	7.62%	1.197%
14	9.014	2234	2252	2256	rBV3	233283	466753	2.38%	0.375%
15	9.236	2312	2321	2334	rVB	121859	242123	1.24%	0.194%
16	9.332	2335	2351	2364	rBV2	97496	235993	1.20%	0.189%
17	9.487	2388	2399	2418	rVB2	288696	584671	2.98%	0.469%
18	9.586	2419	2430	2433	rBV2	160129	262171	1.34%	0.210%
19	9.622	2434	2441	2456	rVB2	326742	718629	3.67%	0.577%
20	9.837	2488	2508	2528	rBV5	110083	478927	2.44%	0.384%
21	10.059	2564	2577	2589	rBV2	900622	1841184	9.40%	1.478%
22	10.123	2590	2597	2608	rVB	302960	543123	2.77%	0.436%
23	10.230	2622	2630	2635	rBV4	113762	203613	1.04%	0.163%
24	10.493	2700	2712	2728	rVB5	431986	946245	4.83%	0.759%
25	11.184	2908	2927	2945	rVB9	68312	218623	1.12%	0.175%
26	11.381	2977	2988	3005	rBV	675426	1243018	6.35%	0.998%
27	11.483	3007	3020	3038	rVV	8866977	15606717	79.67%	12.526%
28	11.599	3043	3056	3076	rVB	3229459	5630523	28.74%	4.519%
29	11.921	3141	3156	3174	rVB	1213993	2088950	10.66%	1.677%
30	12.223	3239	3250	3261	rBV	610572	976969	4.99%	0.784%
31	12.377	3283	3298	3315	rBV	513494	914879	4.67%	0.734%
32	12.564	3344	3356	3370	rVB	1302043	2137084	10.91%	1.715%
33	12.664	3373	3387	3395	rVV	877729	1424734	7.27%	1.144%
34	12.709	3395	3401	3412	rVB	180219	279406	1.43%	0.224%

LSC Area Percent Report

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Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

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\82N070820W.M
Title : SW846 8260

35	12.866	3437	3450	3465	rBV	1788942	2927706	14.94%	2.350%
36	13.014	3482	3496	3509	rBV	11792392	19590088	100.00%	15.723%
37	13.136	3521	3534	3547	rVB3	202487	498107	2.54%	0.400%
38	13.348	3581	3600	3612	rBV	3504890	6590256	33.64%	5.289%
39	13.519	3631	3653	3662	rBV	5575577	10747100	54.86%	8.626%
40	13.567	3662	3668	3684	rVB2	1040825	1916599	9.78%	1.538%
41	13.647	3684	3693	3699	rBV	162463	242086	1.24%	0.194%
42	13.699	3699	3709	3723	rVB2	923393	1864490	9.52%	1.496%
43	13.779	3723	3734	3740	rBV2	1124531	1986631	10.14%	1.594%
44	13.863	3751	3760	3770	rVB	1561447	2326981	11.88%	1.868%
45	13.921	3770	3778	3784	rBV	317242	469950	2.40%	0.377%
46	13.969	3784	3793	3807	rVB	1436589	2342813	11.96%	1.880%
47	14.104	3825	3835	3844	rVB2	266565	408380	2.08%	0.328%
48	14.184	3844	3860	3866	rBV	1418547	2323776	11.86%	1.865%
49	14.223	3866	3872	3881	rVV	1623376	2366952	12.08%	1.900%
50	14.268	3881	3886	3893	rVB2	191863	239409	1.22%	0.192%
51	14.451	3934	3943	3952	rVV	1150003	1828024	9.33%	1.467%
52	14.496	3952	3957	3965	rVB2	258443	368963	1.88%	0.296%
53	14.567	3965	3979	3990	rBV3	1751354	3550632	18.12%	2.850%
54	14.628	3990	3998	4010	rVB4	371895	627811	3.20%	0.504%
55	14.715	4014	4025	4034	rBV	468073	779691	3.98%	0.626%
56	14.824	4040	4059	4067	rBV4	813172	2019793	10.31%	1.621%
57	14.930	4079	4092	4111	rVB3	715964	1602067	8.18%	1.286%
58	15.104	4125	4146	4158	rBV	3136773	5543513	28.30%	4.449%
59	15.207	4167	4178	4188	rBV3	318407	632915	3.23%	0.508%
60	15.261	4188	4195	4221	rVB	179936	368133	1.88%	0.295%
61	15.387	4221	4234	4247	rBV	390301	724879	3.70%	0.582%
62	15.544	4266	4283	4302	rBV	242105	687808	3.51%	0.552%
63	15.667	4311	4321	4330	rBV	126639	235478	1.20%	0.189%
64	15.734	4331	4342	4354	rVB4	155463	339214	1.73%	0.272%
65	16.123	4452	4463	4486	rVB	245448	532189	2.72%	0.427%
66	16.313	4507	4522	4539	rBV	350115	764352	3.90%	0.613%

Sum of corrected areas: 124592850

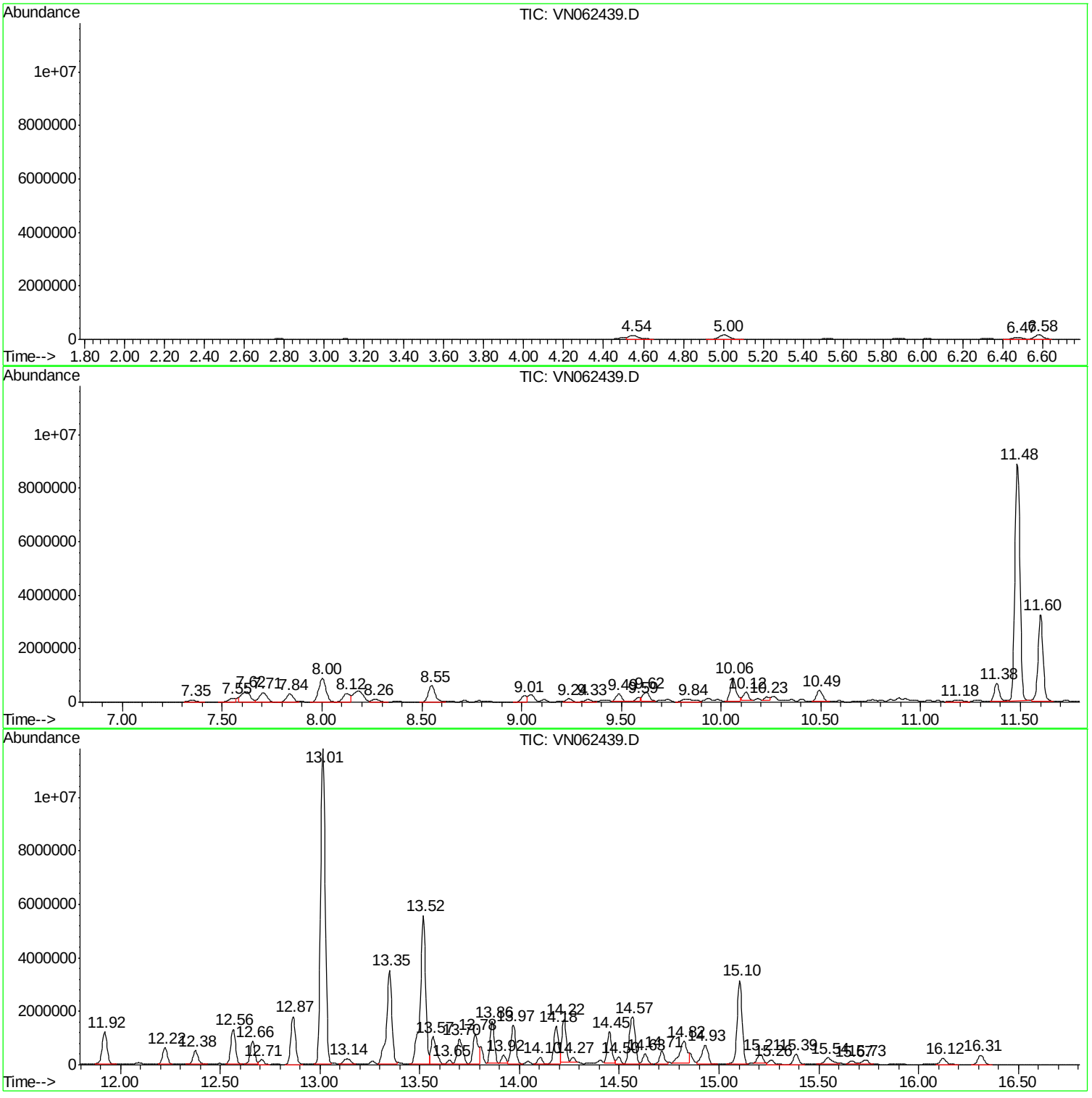
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.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_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

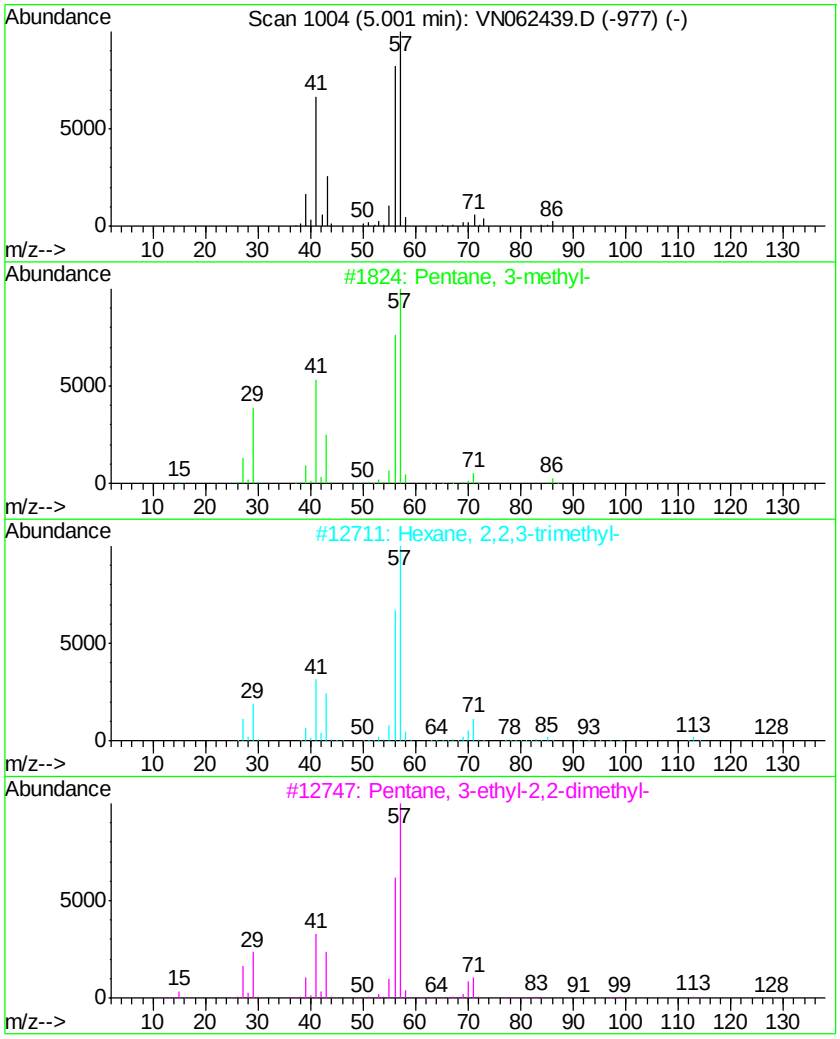
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

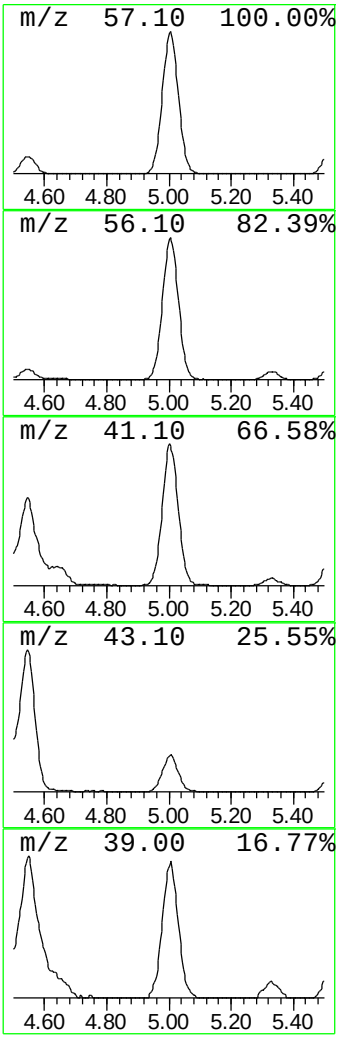
****

Peak Number 1 Pentane, 3-methyl- Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
5.00	24.43 ug/l	599884	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 3-methyl-	86	C6H14	000096-14-0	91
2	Hexane, 2,2,3-trimethyl-	128	C9H20	016747-25-4	64
3	Pentane, 3-ethyl-2,2-dimethyl-	128	C9H20	016747-32-3	56
4	3,4-Dimethyldihydrofuran-2,5-dione	128	C6H8O3	007475-92-5	53
5	Pentane, 2,2,3-trimethyl-	114	C8H18	000564-02-3	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

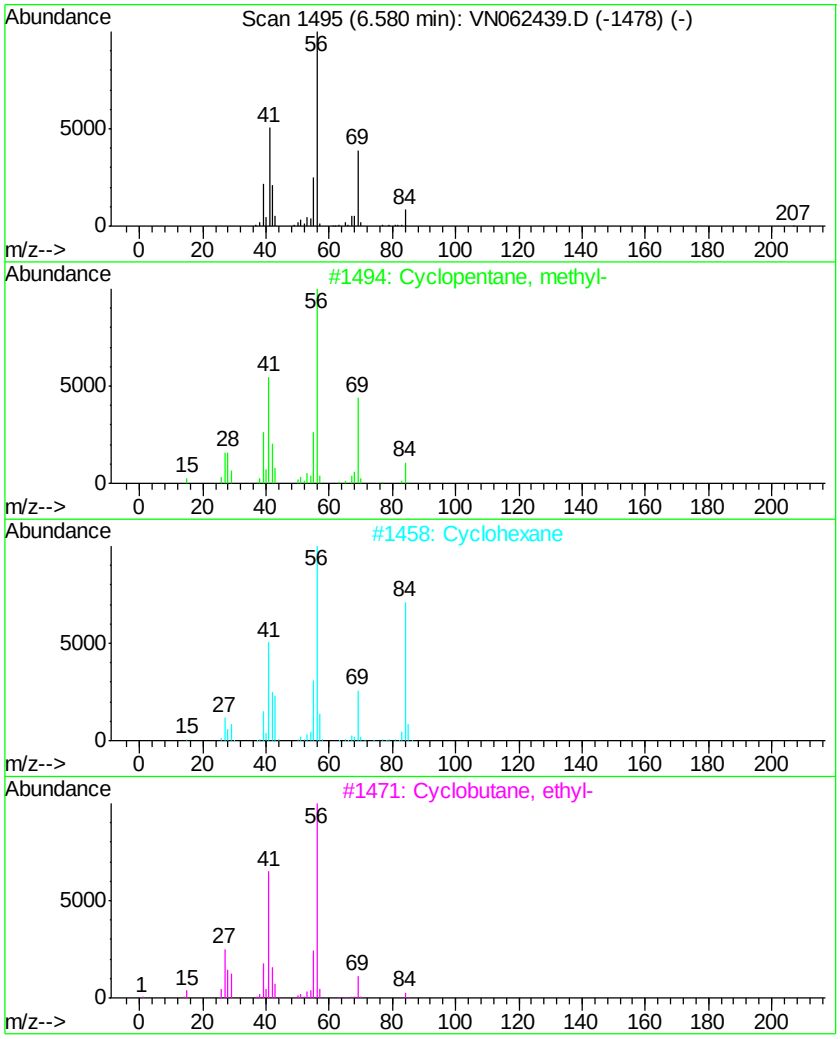
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

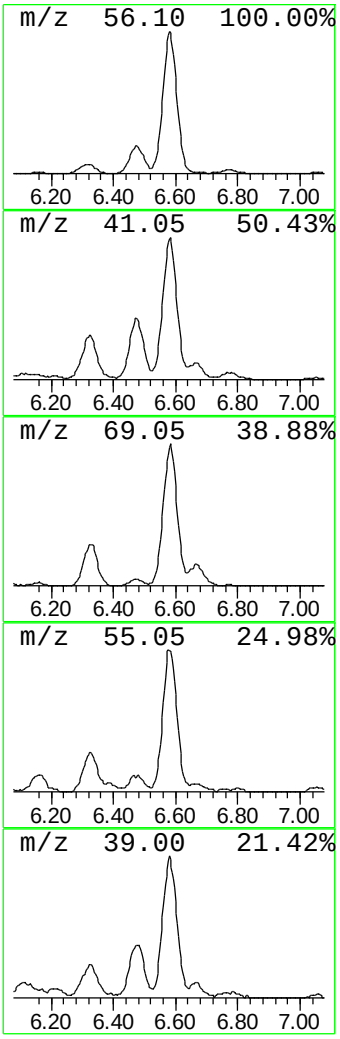
****

Peak Number 2 Cyclopentane, methyl- Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
6.58	19.88 ug/l	488238	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, methyl-	84	C6H12	000096-37-7	90
2	Cyclohexane	84	C6H12	000110-82-7	83
3	Cyclobutane, ethyl-	84	C6H12	004806-61-5	72
4	1H-Tetrazole, 5-methyl-	84	C2H4N4	004076-36-2	64
5	Cyclopentane, 1,1-dimethyl-	98	C7H14	001638-26-2	39





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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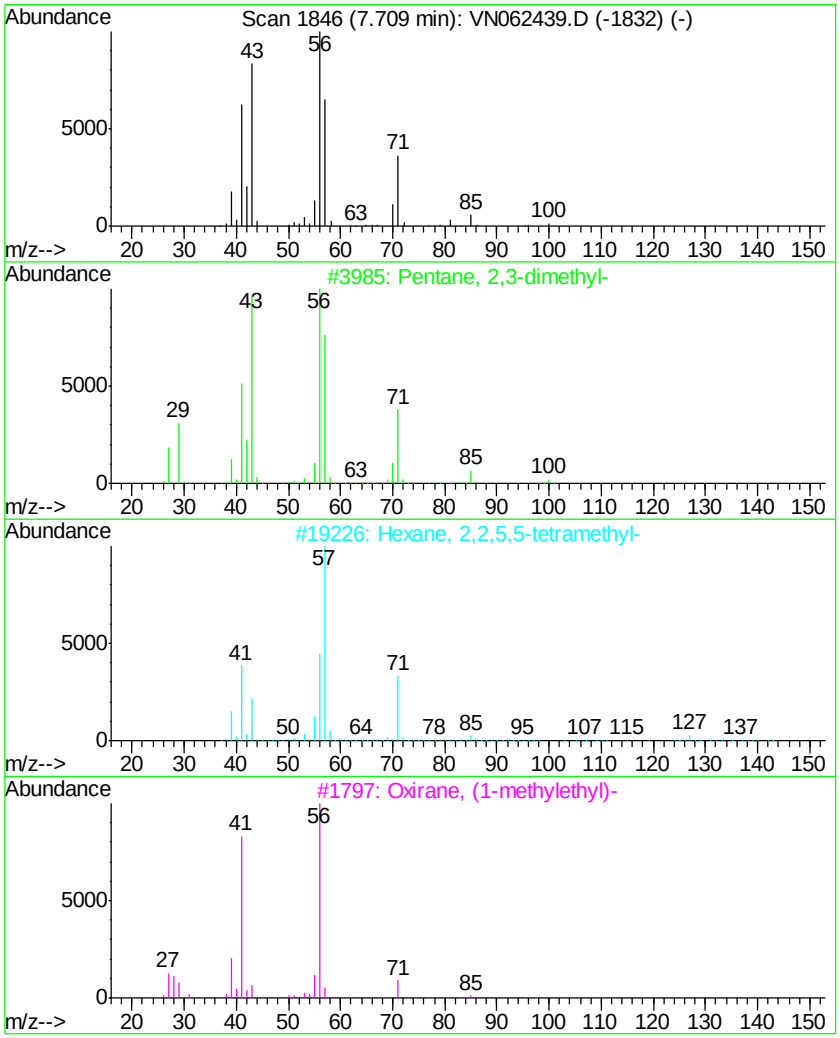
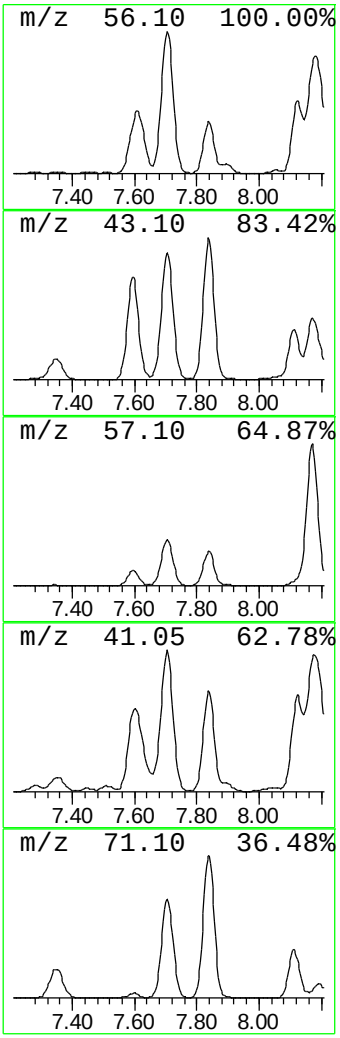
Peak Number 3 Pentane, 2,3-dimethyl- Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.71	39.50 ug/l	969916	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3-dimethyl-	100	C7H16	000565-59-3	91
2	Hexane, 2,2,5,5-tetramethyl-	142	C10H22	001071-81-4	45
3	Oxirane, (1-methylethyl)-	86	C5H10O	001438-14-8	43
4	Butane, 2,2,3-trimethyl-	100	C7H16	000464-06-2	40
5	1-Pentene, 4-methyl-	84	C6H12	000691-37-2	38





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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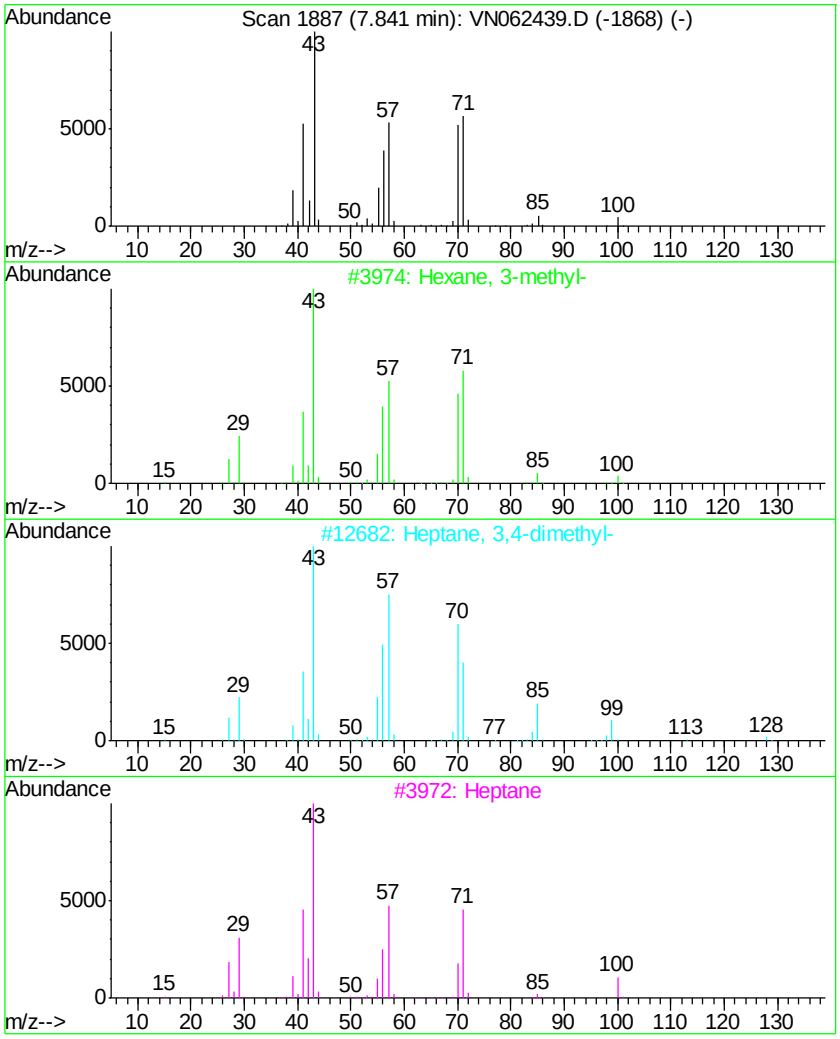
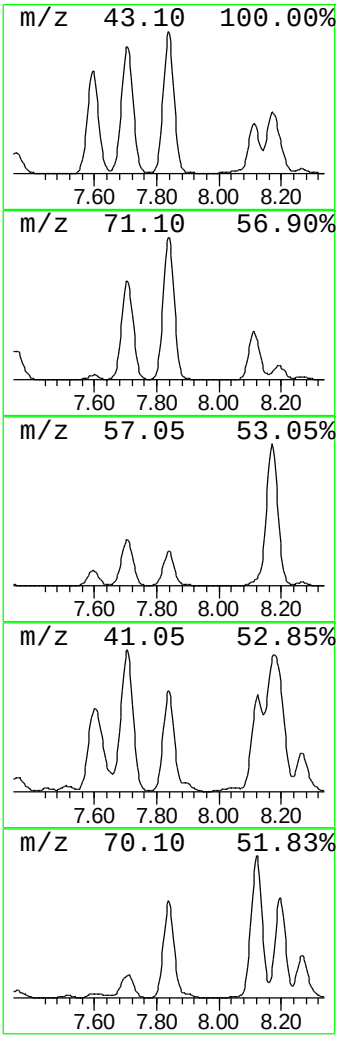
Peak Number 4 Hexane, 3-methyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
7.84	30.61 ug/l	751618	Pentafluorobenzene	7.63

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Hexane, 3-methyl-	100	C7H16	000589-34-4	94
2	Heptane, 3,4-dimethyl-	128	C9H20	000922-28-1	64
3	Heptane	100	C7H16	000142-82-5	64
4	Heptane, 4-methyl-	114	C8H18	000589-53-7	59
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	53





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

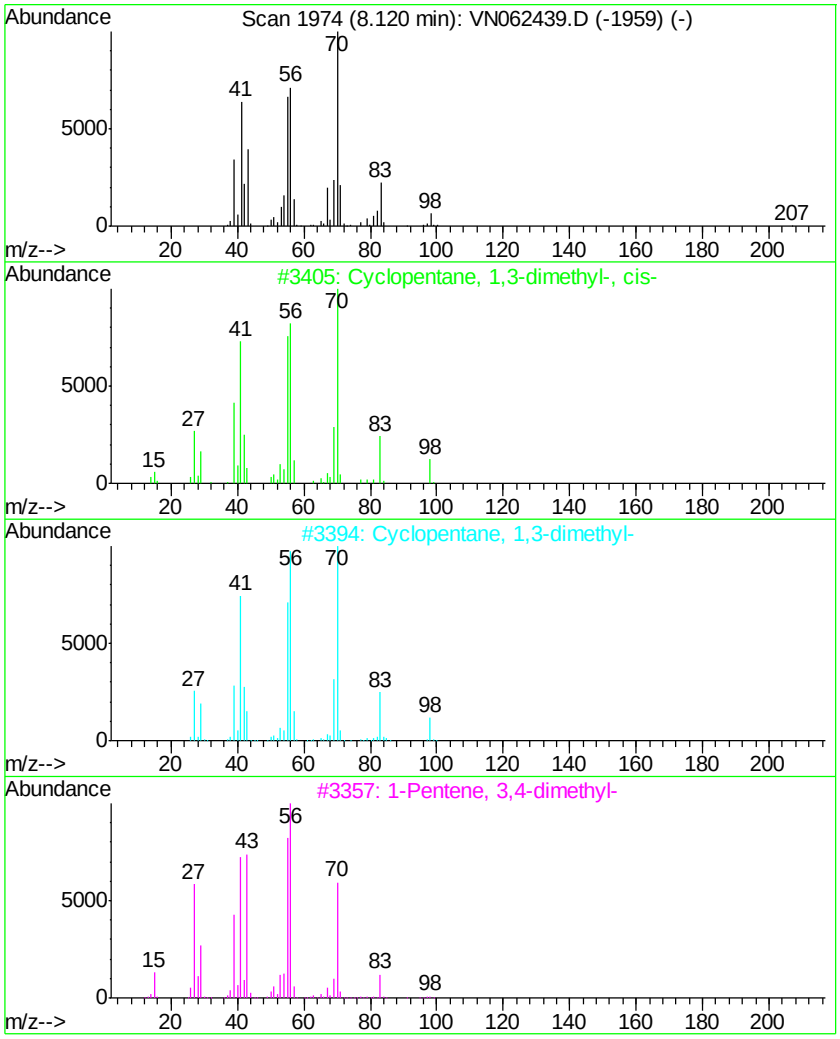
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

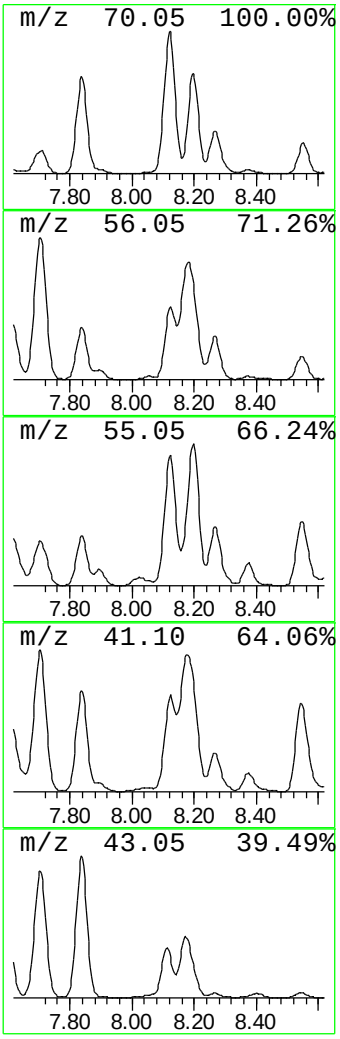
****

Peak Number 5 Cyclopentane, 1,3-dimethyl-... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
8.12	26.22 ug/l	782424	1,4-Difluorobenzene	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentane, 1,3-dimethyl-, cis-	98	C7H14	002532-58-3	72
2	Cyclopentane, 1,3-dimethyl-	98	C7H14	002453-00-1	64
3	1-Pentene, 3,4-dimethyl-	98	C7H14	007385-78-6	53
4	1-Nonanol	144	C9H20O	000143-08-8	47
5	Isopropylcyclobutane	98	C7H14	000872-56-0	45





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

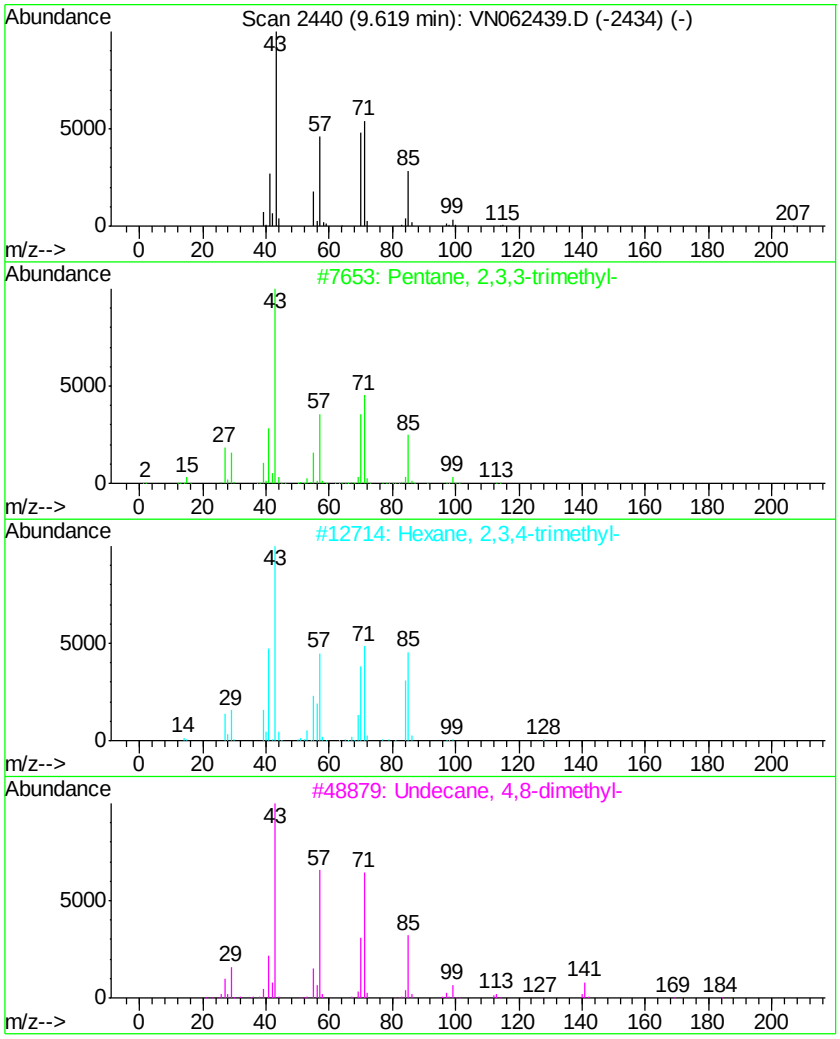
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

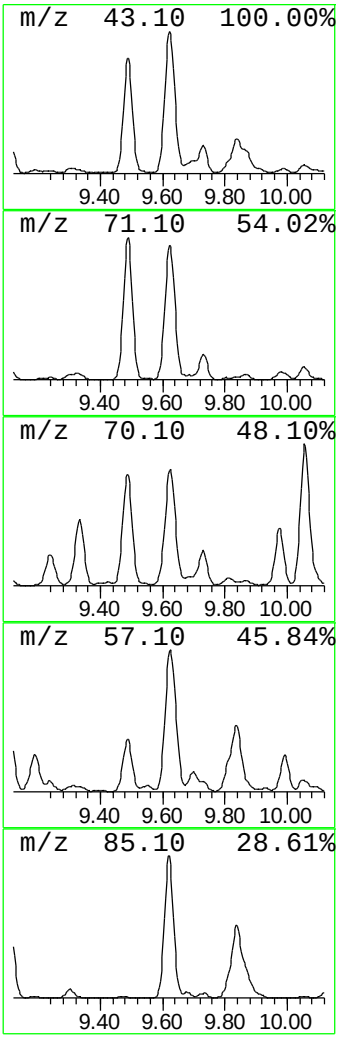
****

Peak Number 6 Pentane, 2,3,3-trimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
9.62	24.08 ug/l	718629	1,4-Difluorobenzene	8.55

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Pentane, 2,3,3-trimethyl-	114	C8H18	000560-21-4	90
2	Hexane, 2,3,4-trimethyl-	128	C9H20	000921-47-1	78
3	Undecane, 4,8-dimethyl-	184	C13H28	017301-33-6	72
4	Hexane, 3-methyl-	100	C7H16	000589-34-4	64
5	Pentane, 2,3,4-trimethyl-	114	C8H18	000565-75-3	64





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

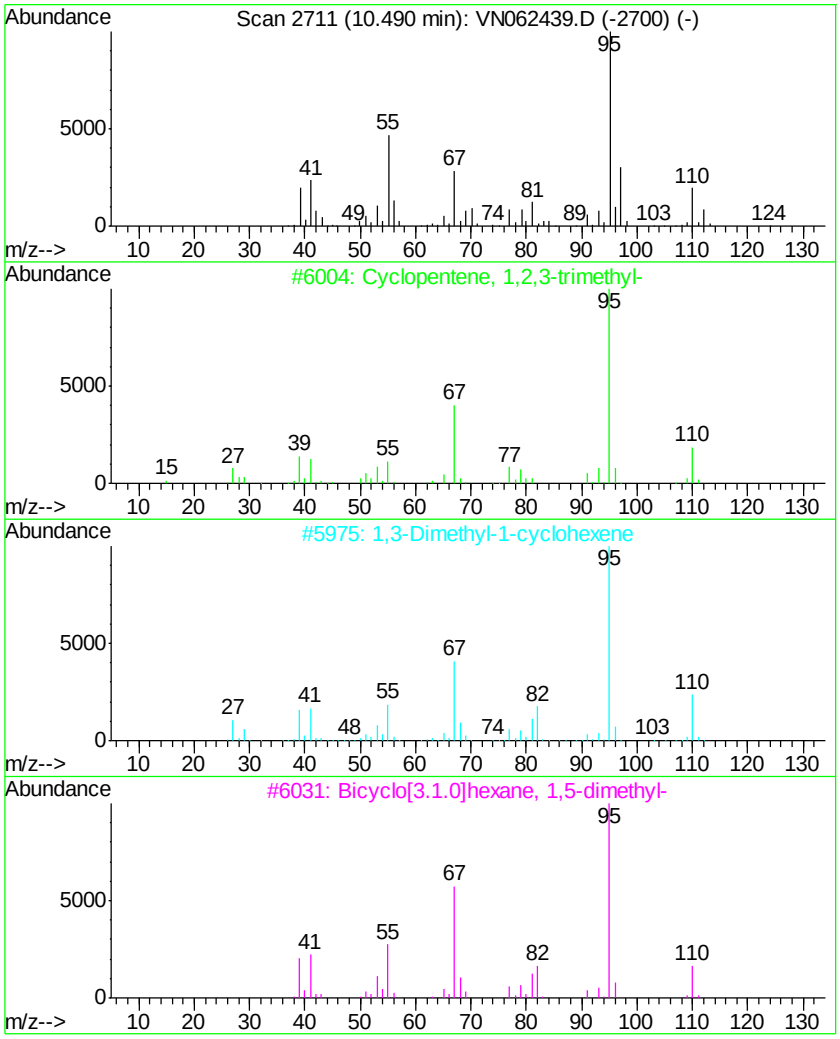
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

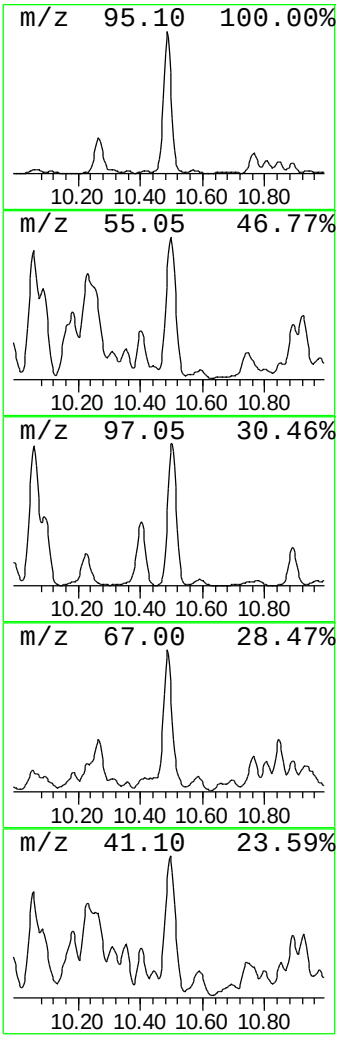
****

Peak Number 7 Cyclopentene, 1,2,3-trimethyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.49	38.06 ug/l	946245	Chlorobenzene-d5	11.38

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Cyclopentene, 1,2,3-trimethyl-	110	C8H14	000473-91-6	90
2	1,3-Dimethyl-1-cyclohexene	110	C8H14	002808-76-6	58
3	Bicyclo[3.1.0]hexane, 1,5-dimethyl-	110	C8H14	1000142-17-5	55
4	Cyclohexene, 3,5-dimethyl-	110	C8H14	000823-17-6	49
5	1,4-Hexadiene, 2,3-dimethyl-	110	C8H14	018669-52-8	46





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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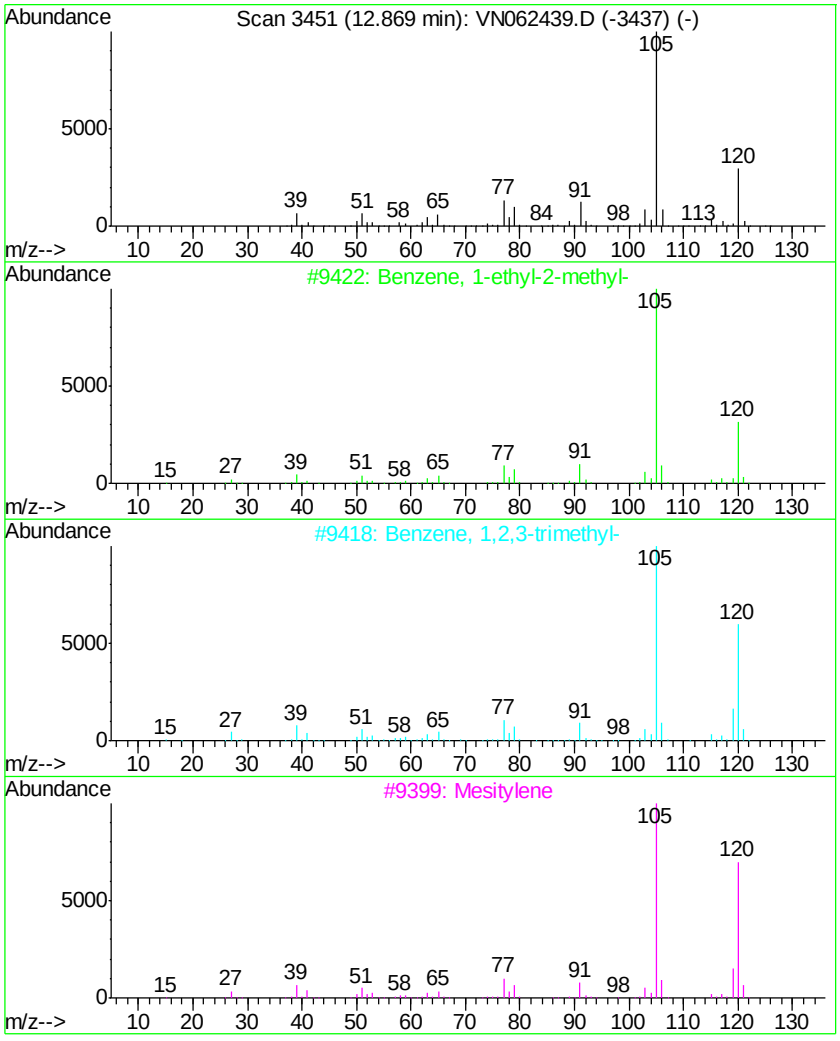
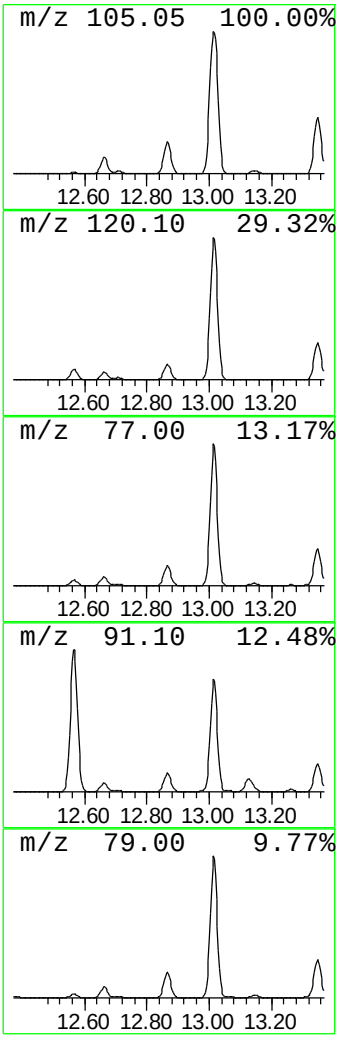
Peak Number 8 Benzene, 1-ethyl-2-methyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
12.87	22.21 ug/l	2927710	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-ethyl-2-methyl-	120	C9H12	000611-14-3	94
2	Benzene, 1,2,3-trimethyl-	120	C9H12	000526-73-8	91
3	Mesitylene	120	C9H12	000108-67-8	91
4	Benzene, 1-ethyl-3-methyl-	120	C9H12	000620-14-4	91
5	Benzene, 1-ethyl-4-methyl-	120	C9H12	000622-96-8	91





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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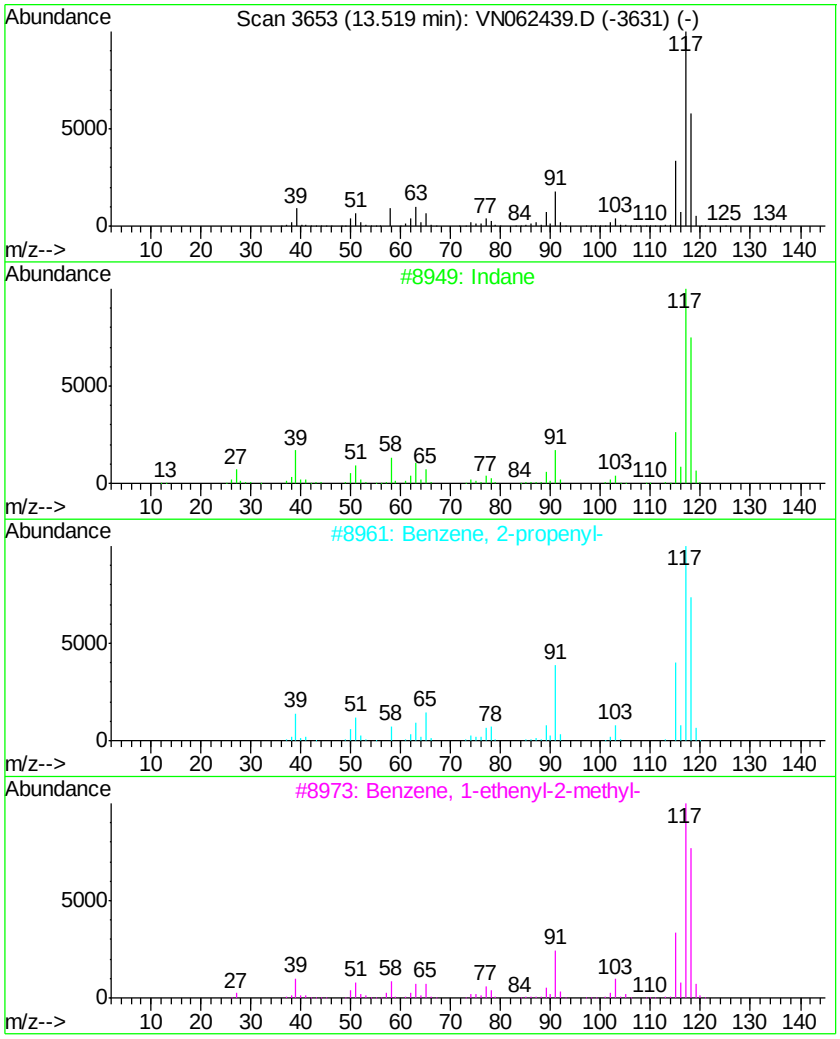
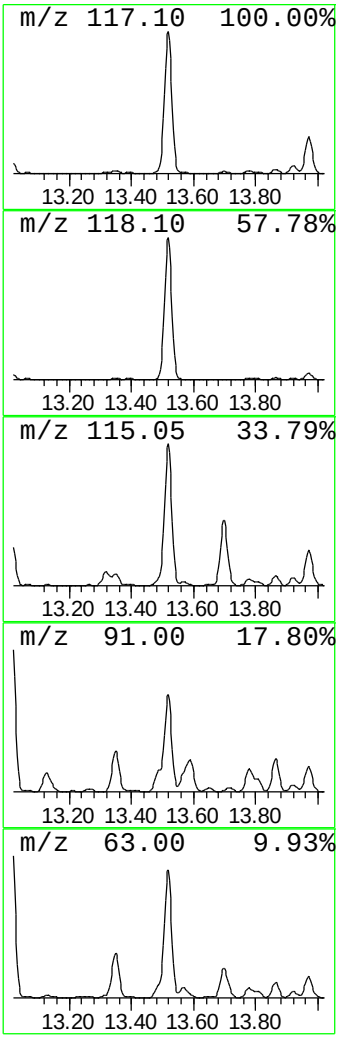
Peak Number 9 Indane Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.52	81.54 ug/l	10747100	1,4-Dichlorobenzene-d4	13.32

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Indane	118	C9H10	000496-11-7	96
2	Benzene, 2-propenyl-	118	C9H10	000300-57-2	81
3	Benzene, 1-ethenyl-2-methyl-	118	C9H10	000611-15-4	76
4	Benzene, cyclopropyl-	118	C9H10	000873-49-4	72
5	Deltacyclene	118	C9H10	007785-10-6	68





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.M
Quant Title : SW846 8260

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

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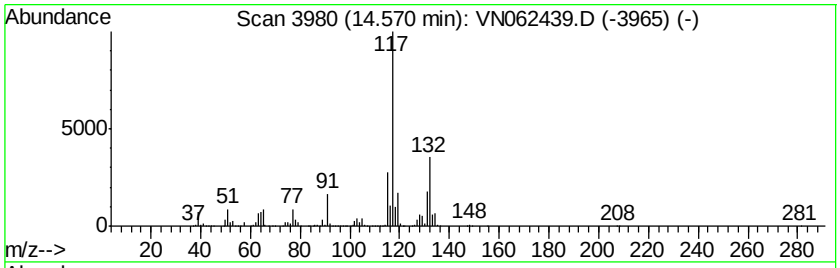
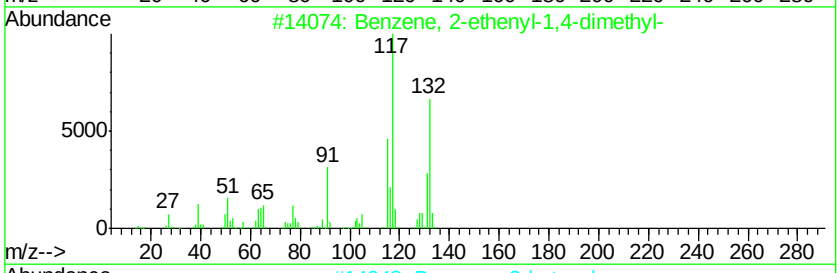
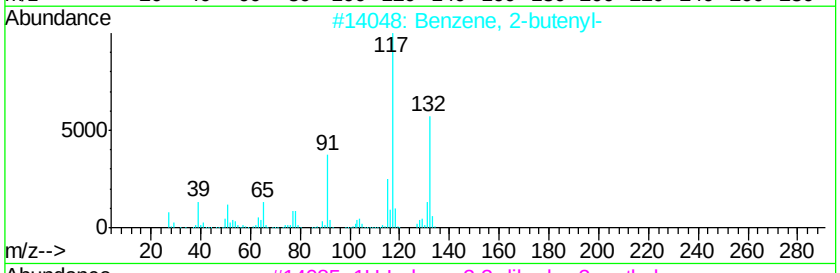
Peak Number 10 Benzene, 2-ethenyl-1,4-dime... Concentration Rank 5

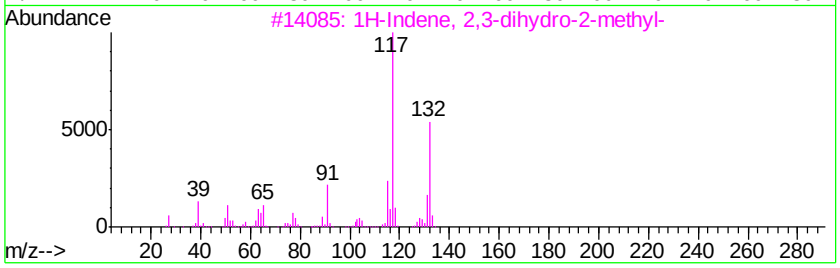
R.T.	EstConc	Area	Relative to ISTD	R.T.
14.57	26.94 ug/l	3550630	1,4-Dichlorobenzene-d4	13.32

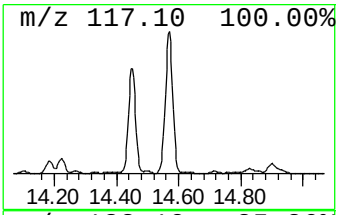
Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 2-ethenyl-1,4-dimethyl-	132	C10H12	002039-89-6	95
2	Benzene, 2-butenyl-	132	C10H12	001560-06-1	87
3	1H-Indene, 2,3-dihydro-2-methyl-	132	C10H12	000824-63-5	87
4	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	83
5	3-Phenylbut-1-ene	132	C10H12	000934-10-1	81

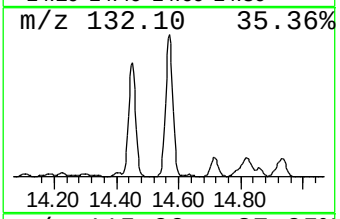


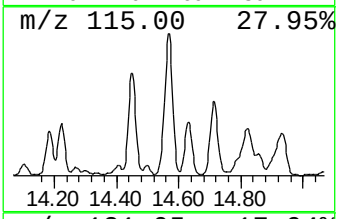


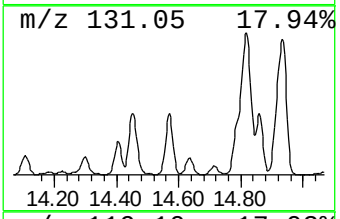


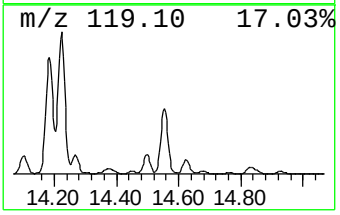












Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN070820\
Data File : VN062439.D
Acq On : 9 Jul 2020 5:11
Operator : JC/MD
Sample : L3215-18
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 49 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : MW-3

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N070820W.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
Pentane, 3-methyl-	5.00	24.4	ug/l	599884	1	7.63	1227860	50.0
Cyclopentane, met...	6.58	19.9	ug/l	488238	1	7.63	1227860	50.0
Pentane, 2,3-dime...	7.71	39.5	ug/l	969916	1	7.63	1227860	50.0
Hexane, 3-methyl-	7.84	30.6	ug/l	751618	1	7.63	1227860	50.0
Cyclopentane, 1,3...	8.12	26.2	ug/l	782424	2	8.55	1491910	50.0
Pentane, 2,3,3-tr...	9.62	24.1	ug/l	718629	2	8.55	1491910	50.0
Cyclopentene, 1,2...	10.49	38.1	ug/l	946245	3	11.38	1243020	50.0
Benzene, 1-ethyl-...	12.87	22.2	ug/l	2927710	4	13.32	6590260	50.0
Indane	13.52	81.5	ug/l	10747100	4	13.32	6590260	50.0
Benzene, 2-etheny...	14.57	26.9	ug/l	3550630	4	13.32	6590260	50.0