LSC Area Percent Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

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\82N031219W.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	7.587	1782	1808	1819	rBV2	951973	2485939	3.91%	0.320%
2	7.661	1819	1831	1847	rVB	1800313	4322345	6.80%	0.556%
3	8.024	1927	1944	1966	rVB2	911434	2111563	3.32%	0.271%
4	8.583	2103	2118	2161	rVB	2428872	5412410	8.51%	0.696%
5	9.075	2246	2271	2283	rBV3	912735	2338191	3.68%	0.301%
6	9.725	2462	2473	2479	rBV	594899	1108724	1.74%	0.143%
7	9.866	2497	2517	2539	rVB3	838811	2376820	3.74%	0.306%
8	10.088	2571	2586	2615	rBV3	4829606	10642589	16.74%	1.368%
9	10.278	2630	2645	2659	rVB6	690346	2021684	3.18%	0.260%
10	10.429	2684	2692	2709	rVB	954325	1769161	2.78%	0.227%
11	10.529	2709	2723	2733	rBV4	669628	1447132	2.28%	0.186%
12	10.715	2771	2781	2791	rBV	657956	1065180	1.68%	0.137%
13	10.818	2804	2813	2825	rBV	437621	825992	1.30%	0.106%
14	10.882	2826	2833	2839	rBV3	446567	730416	1.15%	0.094%
15	10.950	2847	2854	2862	rVB	1321867	1937595	3.05%	0.249%
16	11.005	2862	2871	2881	rVB	2580738	4434198	6.97%	0.570%
17	11.120	2895	2907	2915	rBV4	834168	1620225	2.55%	0.208%
18	11.204	2915	2933	2951	rVB7	2074369	5847162	9.20%	0.752%
19	11.297	2951	2962	2985	rBV2	1712869	3544695	5.58%	0.456%
20	11.406	2985	2996	3015	rBV	3234633	6205545	9.76%	0.798%
21	11.506	3016	3027	3031	rBV2	513863	960790	1.51%	0.124%
22	11.538	3033	3037	3044	rVB	580241	738136	1.16%	0.095%
23	11.590	3044	3053	3062	rVB6	933626	1751911	2.76%	0.225%
24	11.654	3063	3073	3080	rBV2	2009371	3749619	5.90%	0.482%
25	11.918	3142	3155	3169	rBV6	2549113	6371137	10.02%	0.819%
26	12.043	3187	3194	3200	rBV2	1042621	1371343	2.16%	0.176%
27	12.078	3200	3205	3209	rVV4	579299	727483	1.14%	0.094%
28	12.117	3210	3217	3223	rVV	2186756	3675785	5.78%	0.473%
29	12.162	3224	3231	3249	rVV5	3400228	7409135	11.65%	0.952%
30	12.246	3252	3257	3266	rVB	1160318	1727123	2.72%	0.222%
31	12.336	3280	3285	3295	rVB3	1600971	2413263	3.80%	0.310%
32	12.400	3296	3305	3314	rBV	2989336	4966922	7.81%	0.638%
33	12.448	3314	3320	3324	rBV2	648309	886583	1.39%	0.114%
34	12.561	3347	3355	3361	rBV3	3253899	5648118	8.88%	0.726%

LSC Area Percent Report

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Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

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\82N031219W.M
Title : SW846 8260

35	12.644	3372	3381	3390	rBV8	1062663	2196495	3.45%	0.282%
36	12.725	3393	3406	3416	rVV6	3769363	9836331	15.47%	1.264%
37	12.779	3417	3423	3433	rVB3	1833840	2882278	4.53%	0.371%
38	12.866	3443	3450	3457	rVB6	1212102	1748856	2.75%	0.225%
39	12.959	3472	3479	3494	rVB3	3478597	6511224	10.24%	0.837%
40	13.168	3533	3544	3552	rBV2	4126907	7296863	11.48%	0.938%
41	13.236	3554	3565	3574	rVV4	2784825	5612904	8.83%	0.722%
42	13.342	3591	3598	3612	rVB	3234668	5880092	9.25%	0.756%
43	13.509	3644	3650	3656	rBV	4293843	6192931	9.74%	0.796%
44	13.612	3669	3682	3689	rBV2	3934029	9021621	14.19%	1.160%
45	13.664	3689	3698	3708	rVB4	7852052	14201895	22.34%	1.826%
46	13.889	3758	3768	3777	rVB	17696515	27359877	43.04%	3.517%
47	13.943	3778	3785	3791	rVV	2664773	4087176	6.43%	0.525%
48	13.995	3791	3801	3811	rVB3	14054162	24676324	38.81%	3.172%
49	14.107	3832	3836	3846	rBV4	991348	1437812	2.26%	0.185%
50	14.172	3848	3856	3860	rVV6	4749589	7805258	12.28%	1.003%
51	14.207	3862	3867	3882	rVB	12230733	20196745	31.77%	2.596%
52	14.291	3884	3893	3900	rBV2	7611468	12602308	19.82%	1.620%
53	14.329	3900	3905	3911	rVB2	2600036	2979451	4.69%	0.383%
54	14.365	3912	3916	3920	rBV	1253045	1246082	1.96%	0.160%
55	14.477	3942	3951	3961	rBV3	9925746	19027650	29.93%	2.446%
56	14.586	3973	3985	3996	rBV4	28938003	59813378	94.08%	7.689%
57	14.648	3996	4004	4015	rVV2	9725750	17095395	26.89%	2.198%
58	14.815	4049	4056	4058	rBV2	5611166	6842864	10.76%	0.880%
59	14.850	4060	4067	4073	rVV3	16568153	27868921	43.84%	3.582%
60	14.885	4074	4078	4089	rVV2	8190895	14405924	22.66%	1.852%
61	14.956	4091	4100	4112	rVB3	17087211	34689132	54.56%	4.459%
62	15.107	4137	4147	4155	rBV5	3811638	7672439	12.07%	0.986%
63	15.188	4163	4172	4180	rBV	13422912	22685455	35.68%	2.916%
64	15.242	4181	4189	4199	rBV3	8288870	19175578	30.16%	2.465%
65	15.416	4232	4243	4256	rBV2	10858308	22674971	35.67%	2.915%
66	15.573	4276	4292	4296	rBV6	8853490	21182500	33.32%	2.723%
67	15.612	4297	4304	4318	rVB	21057982	37933614	59.67%	4.876%
68	15.699	4324	4331	4340	rVB6	5688301	9620269	15.13%	1.237%
69	15.766	4342	4352	4362	rBV7	6777752	14542876	22.88%	1.869%
70	15.911	4384	4397	4407	rBV3	12657134	26558574	41.78%	3.414%
71	16.101	4441	4456	4465	rBV4	14327132	32674838	51.40%	4.200%

LSC Area Percent Report

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ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Integration Parameters: RTEINT.P
Integrator: RTE
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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\82N031219W.M
Title : SW846 8260

72	16.159	4465	4474	4499	rVB2	27762067	63574968	100.00%	8.172%
73	16.352	4521	4534	4547	rBV	17963561	41472964	65.23%	5.331%

Sum of corrected areas: 777929747

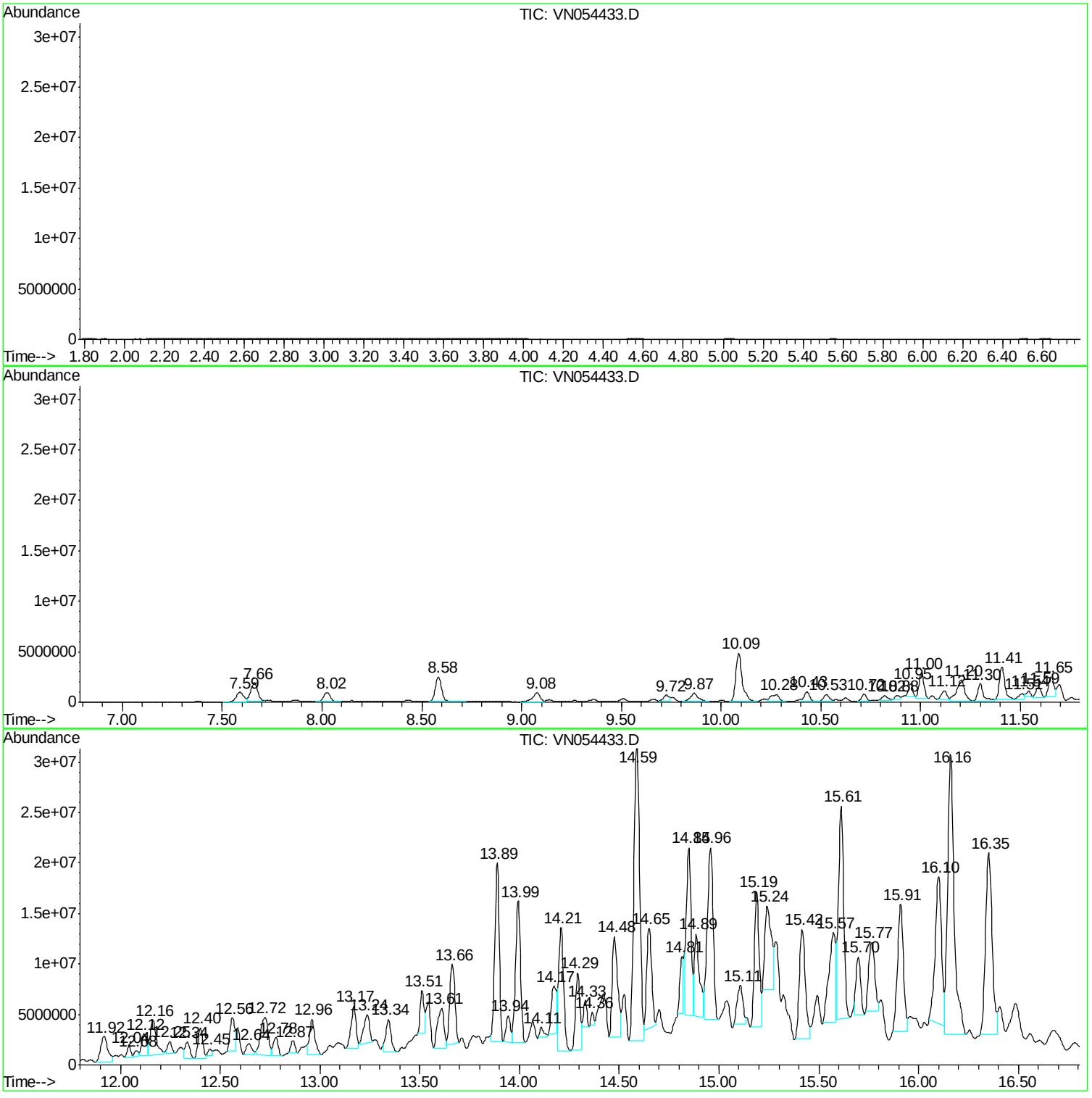
LSC Report - Integrated Chromatogram

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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



Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

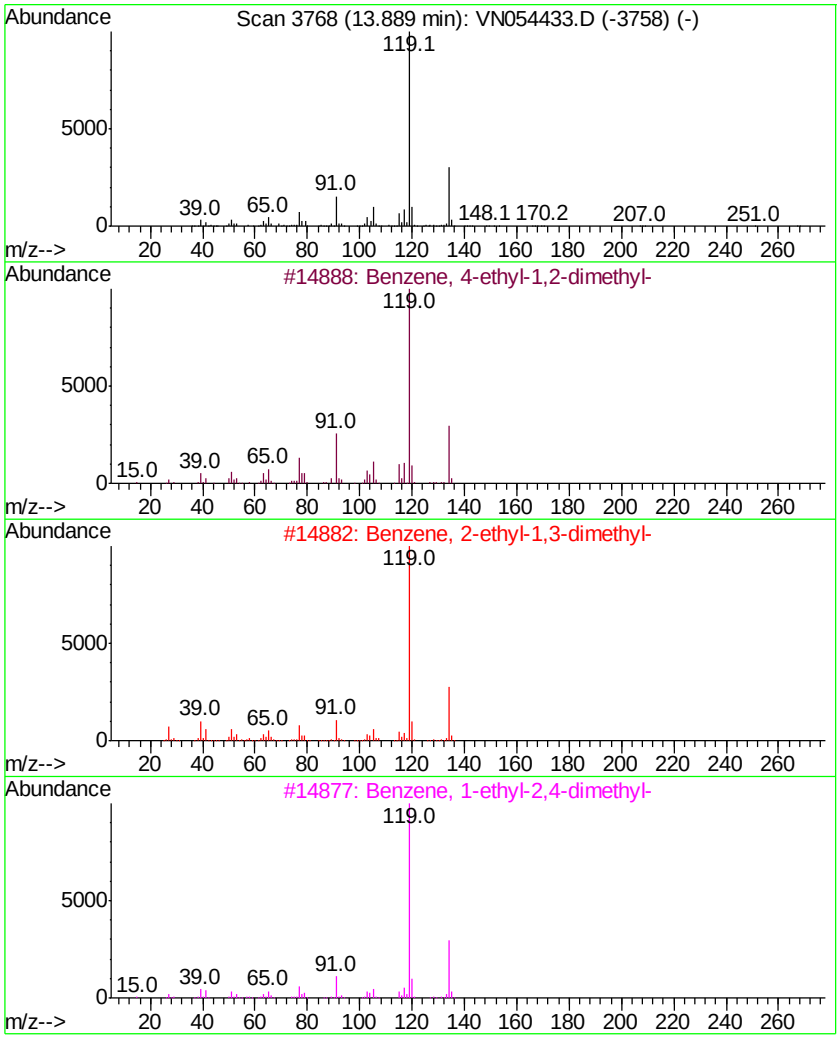
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

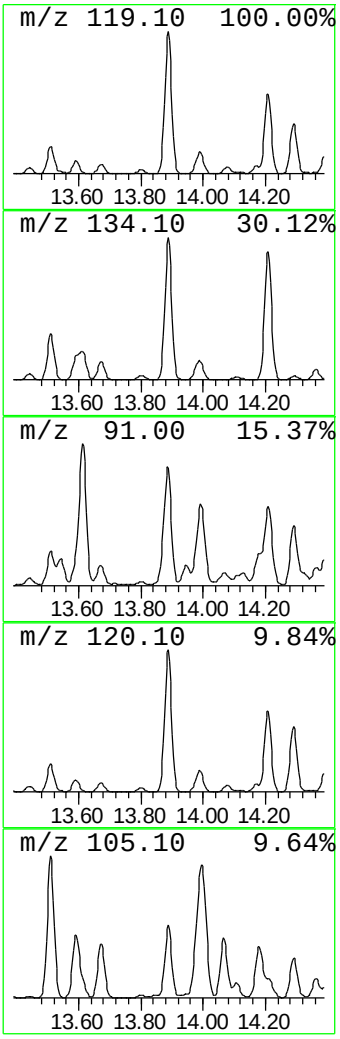
****

Peak Number 1 Benzene, 4-ethyl-1,2-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	232.65 ug/l	27359900	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 4-ethyl-1,2-dimethyl-	134	C10H14	000934-80-5	96
2	Benzene, 2-ethyl-1,3-dimethyl-	134	C10H14	002870-04-4	95
3	Benzene, 1-ethyl-2,4-dimethyl-	134	C10H14	000874-41-9	95
4	o-Cymene	134	C10H14	000527-84-4	95
5	Benzene, 2-ethyl-1,4-dimethyl-	134	C10H14	001758-88-9	94





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

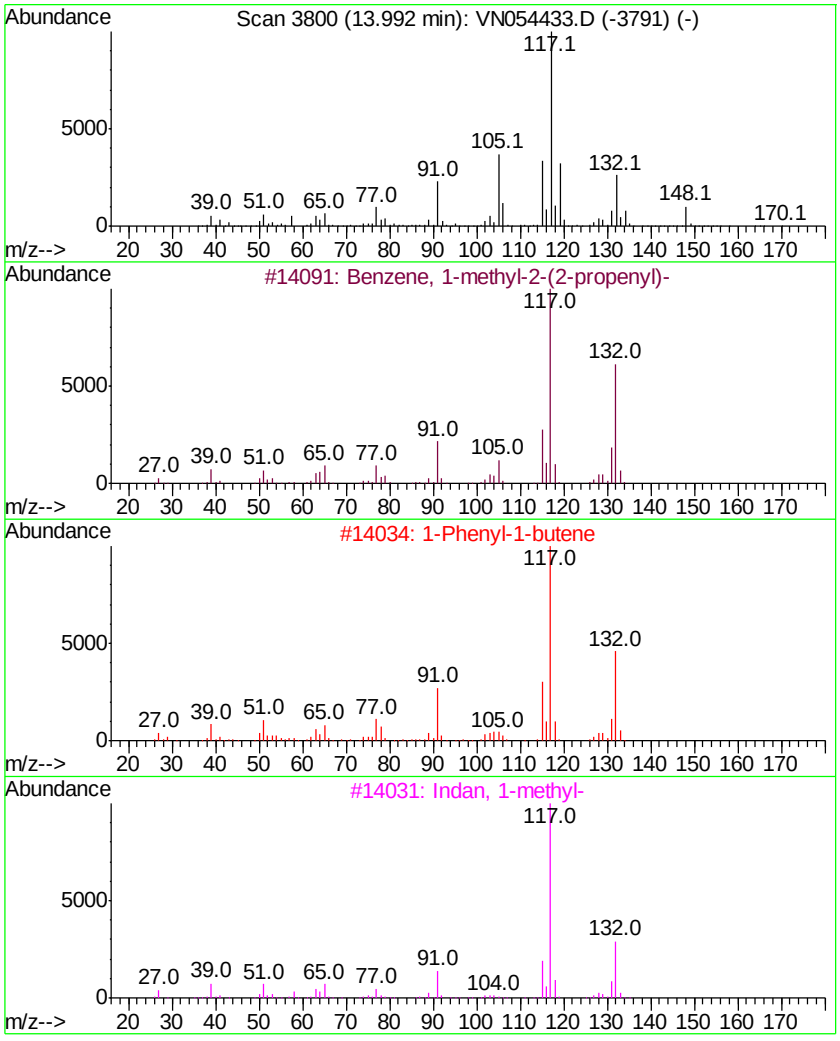
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

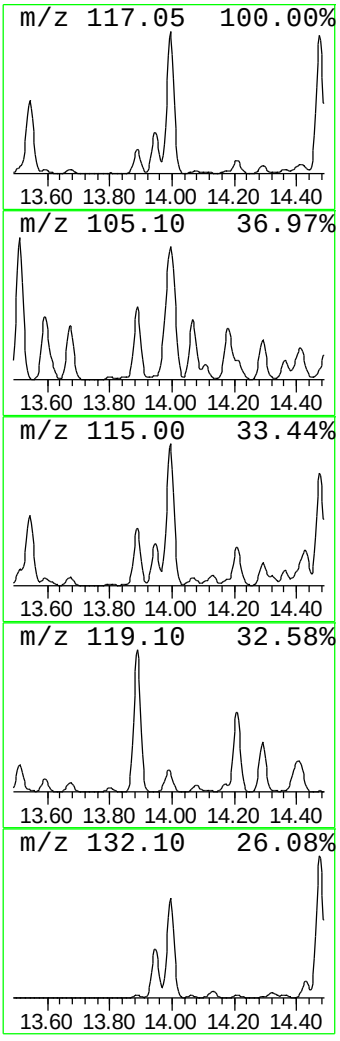
****

Peak Number 2 Benzene, 1-methyl-2-(2-prop... Concentration Rank 10

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.99	209.83 ug/l	24676300	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, 1-methyl-2-(2-propenyl)-	132	C10H12	001587-04-8	76
2	1-Phenyl-1-butene	132	C10H12	000824-90-8	70
3	Indan, 1-methyl-	132	C10H12	000767-58-8	70
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	62
5	Benzene, 2-butenyl-	132	C10H12	001560-06-1	62





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

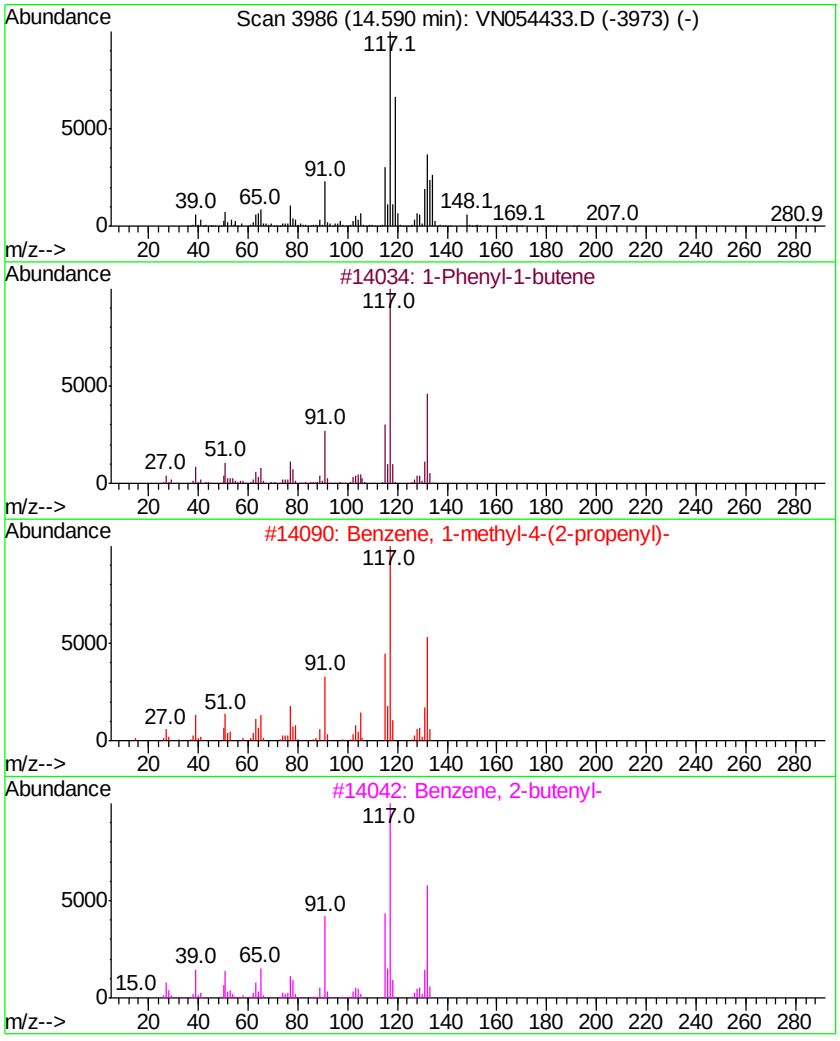
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

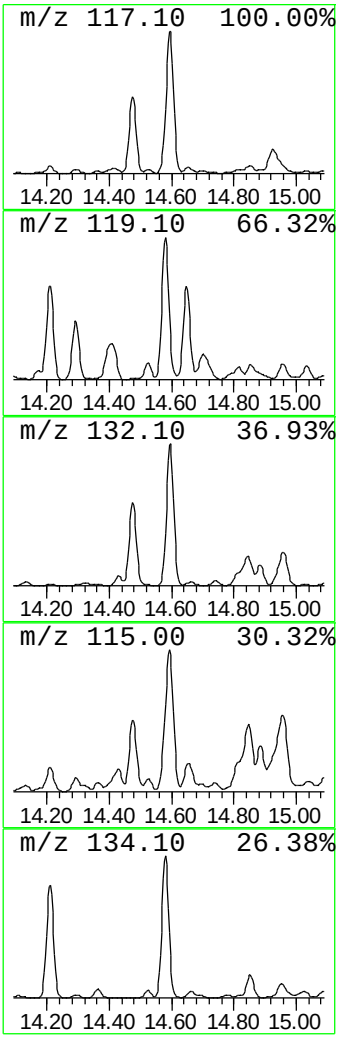
****

Peak Number 3 1-Phenyl-1-butene Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.59	508.61 ug/l	59813400	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	1-Phenyl-1-butene	132	C10H12	000824-90-8	86
2	Benzene, 1-methyl-4-(2-propenyl)-	132	C10H12	003333-13-9	70
3	Benzene, 2-butenyl-	132	C10H12	001560-06-1	64
4	3-Phenylbut-1-ene	132	C10H12	000934-10-1	50
5	Benzene, 2-ethenyl-1,3-dimethyl-	132	C10H12	002039-90-9	50





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
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Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

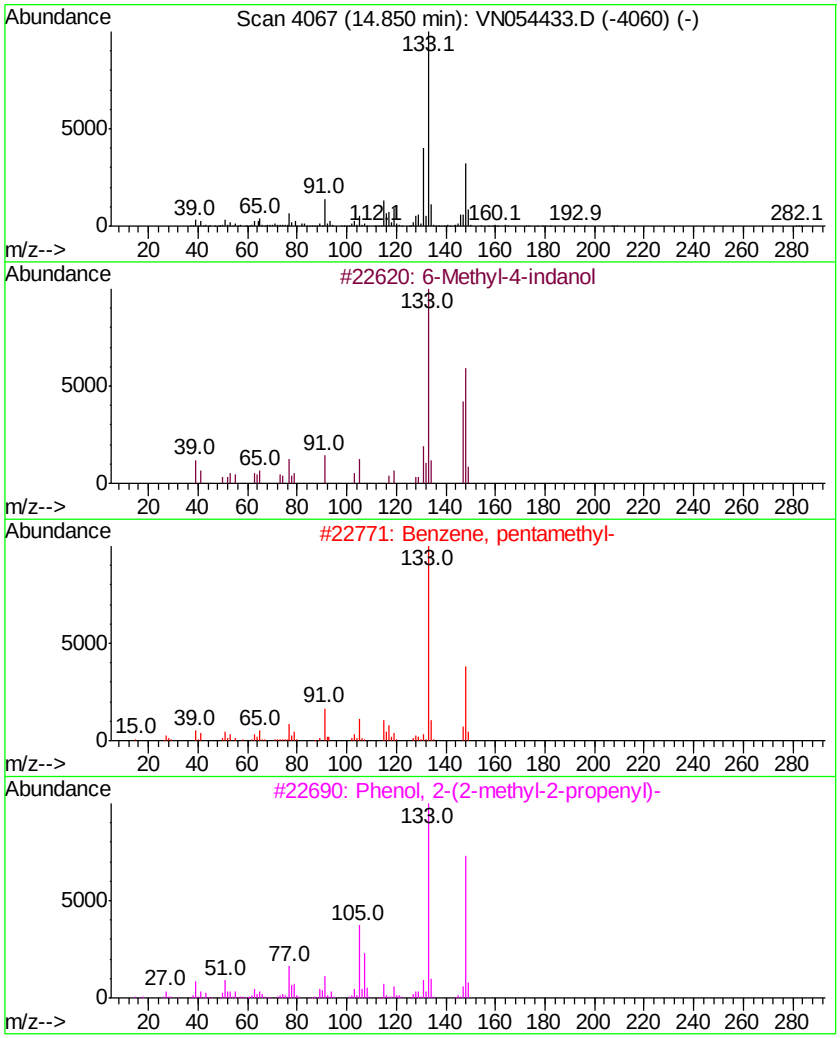
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

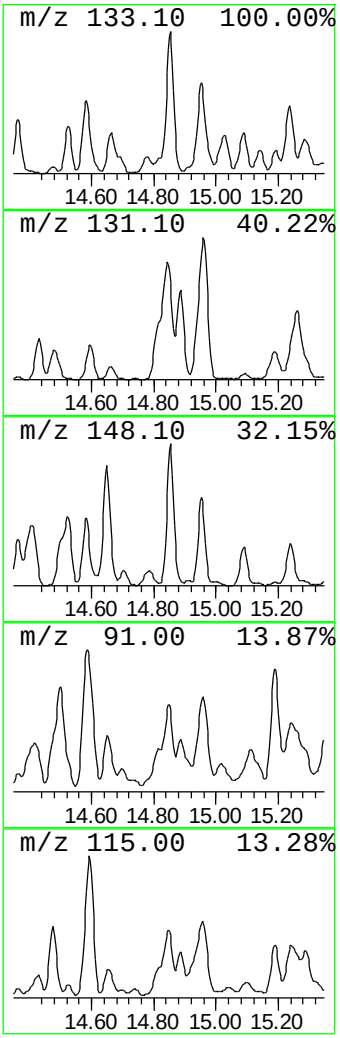
****

Peak Number 4 6-Methyl-4-indanol Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.85	236.98 ug/l	27868900	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	6-Methyl-4-indanol	148	C10H12O	020294-32-0	72
2	Benzene, pentamethyl-	148	C11H16	000700-12-9	58
3	Phenol, 2-(2-methyl-2-propenyl)-	148	C10H12O	020944-88-1	53
4	Pyrazine, 3,5-dimethyl-2-(1-prop...	148	C9H12N2	055138-74-4	52
5	3,4-Dimethylcumene	148	C11H16	1000370-34-1	52





Library Search Compound Report

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ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

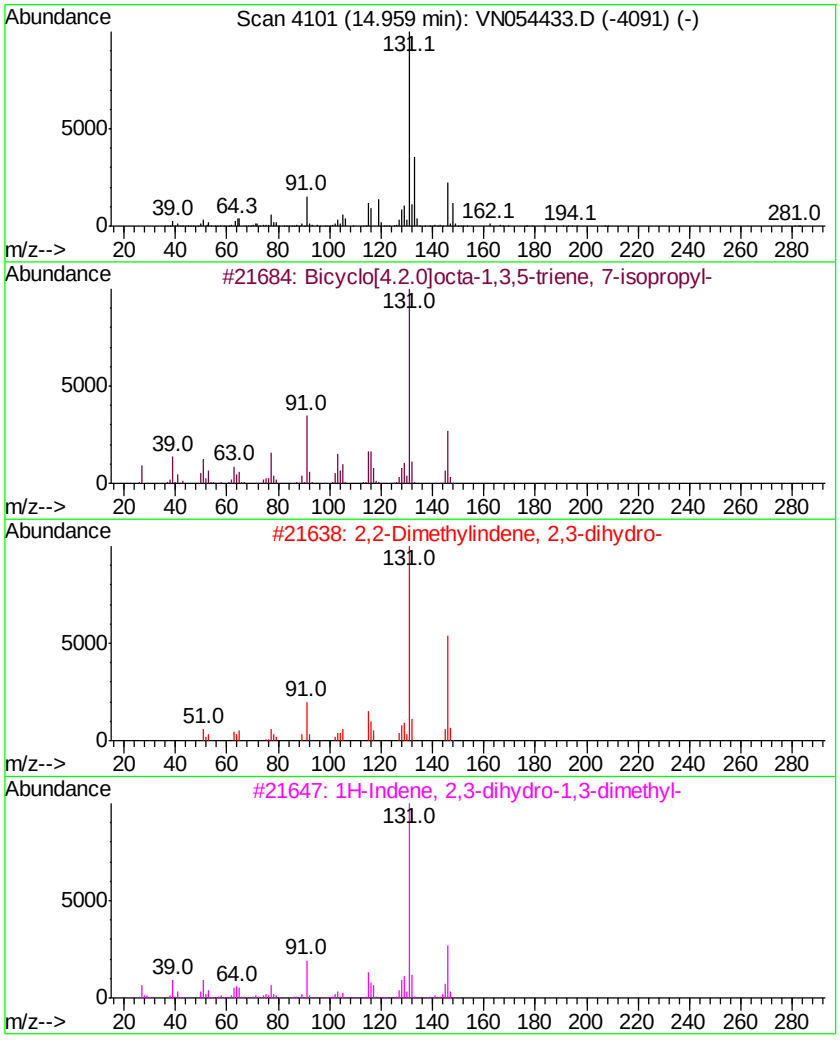
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

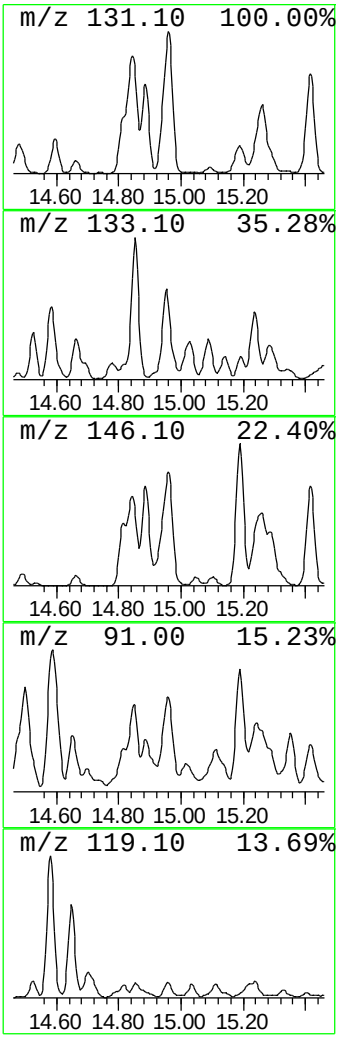
****

Peak Number 5 Bicyclo[4.2.0]octa-1,3,5-tr... Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	294.97 ug/l	34689100	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Bicyclo[4.2.0]octa-1,3,5-triene,...	146	C11H14	027087-54-3	83
2	2,2-Dimethylindene, 2,3-dihydro-	146	C11H14	020836-11-7	64
3	1H-Indene, 2,3-dihydro-1,3-dimet...	146	C11H14	004175-53-5	64
4	1H-Indene, 2,3-dihydro-1,1-dimet...	146	C11H14	004912-92-9	64
5	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001559-81-5	64





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Quant Title : SW846 8260

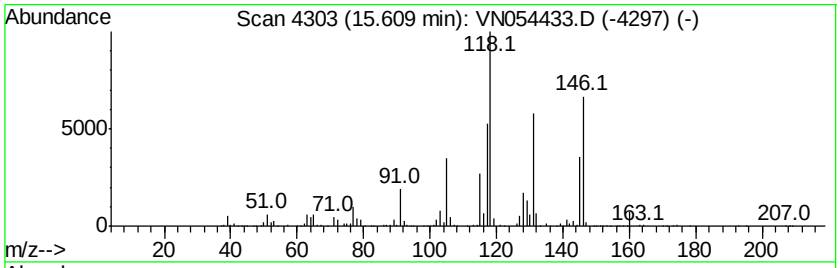
TIC Library : C:\DATABASE\NIST11.L
TIC Integration Parameters: LSCINT.P

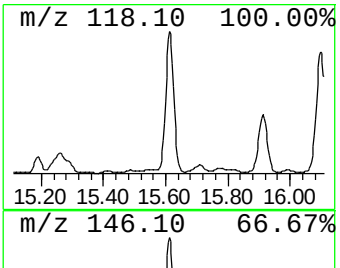
****

Peak Number 6 Benzene, (1-methylenebutyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.61	322.56 ug/l	37933600	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Benzene, (1-methylenebutyl)-	146	C11H14	005676-32-4	87
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	72
3	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	58
4	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	43
5	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	41





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Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

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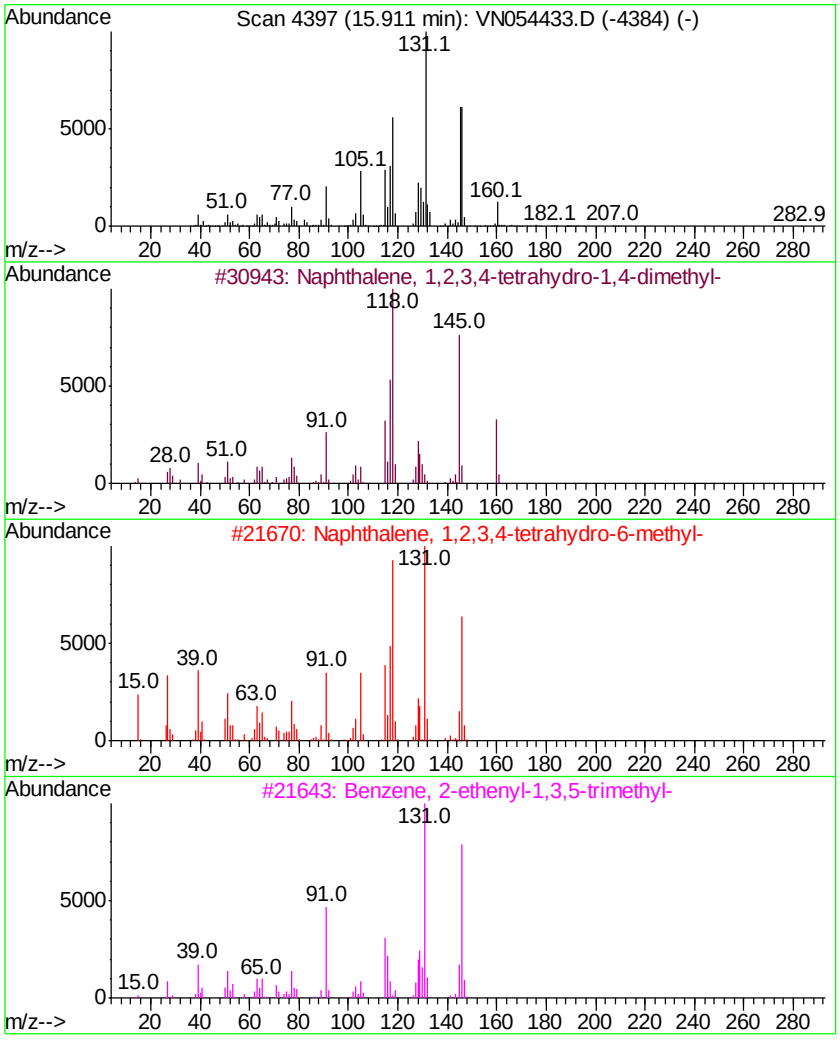
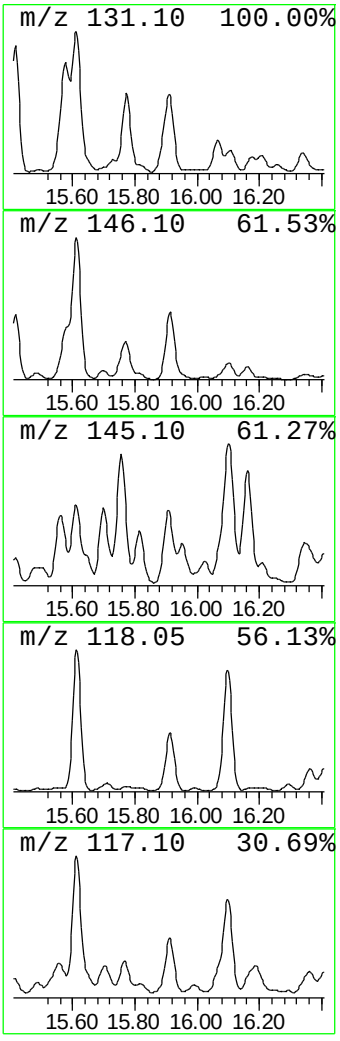
Peak Number 7 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.91	225.84 ug/l	26558600	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	004175-54-6	87
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	86
3	Benzene, 2-ethenyl-1,3,5-trimethyl-	146	C11H14	000769-25-5	83
4	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	81
5	1H-Benzimidazole, 5,6-dimethyl-	146	C9H10N2	000582-60-5	53





Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

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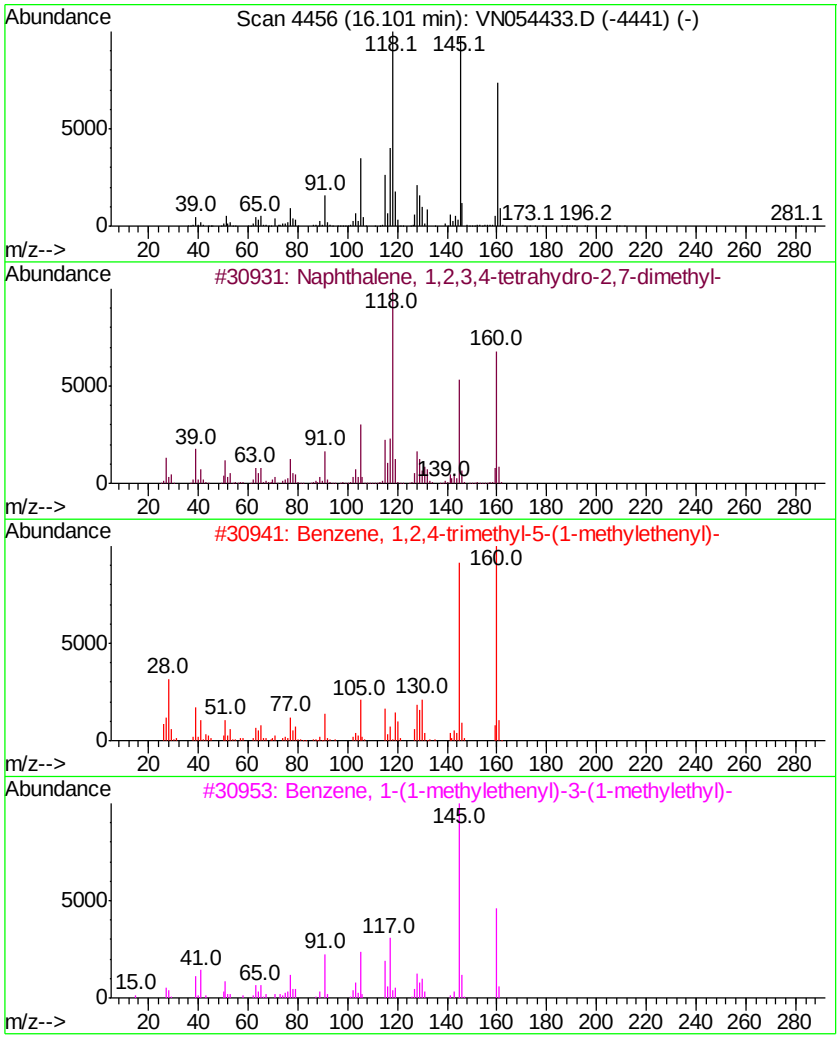
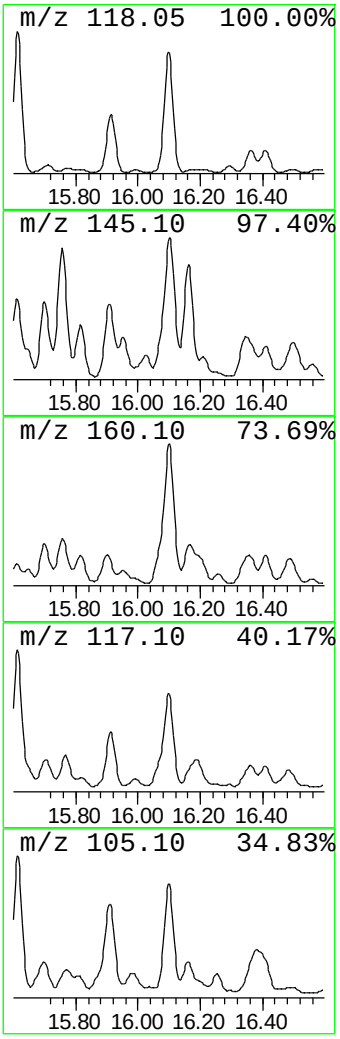
Peak Number 8 Naphthalene, 1,2,3,4-tetrah... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.10	277.84 ug/l	32674800	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	013065-07-1	90
2	Benzene, 1,2,4-trimethyl-5-(1-me...	160	C12H16	054340-84-0	78
3	Benzene, 1-(1-methylethenyl)-3-(...	160	C12H16	001129-29-9	70
4	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	025419-33-4	68
5	Naphthalene, 1,2,3,4-tetrahydro-...	160	C12H16	021693-54-9	64





Library Search Compound Report

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

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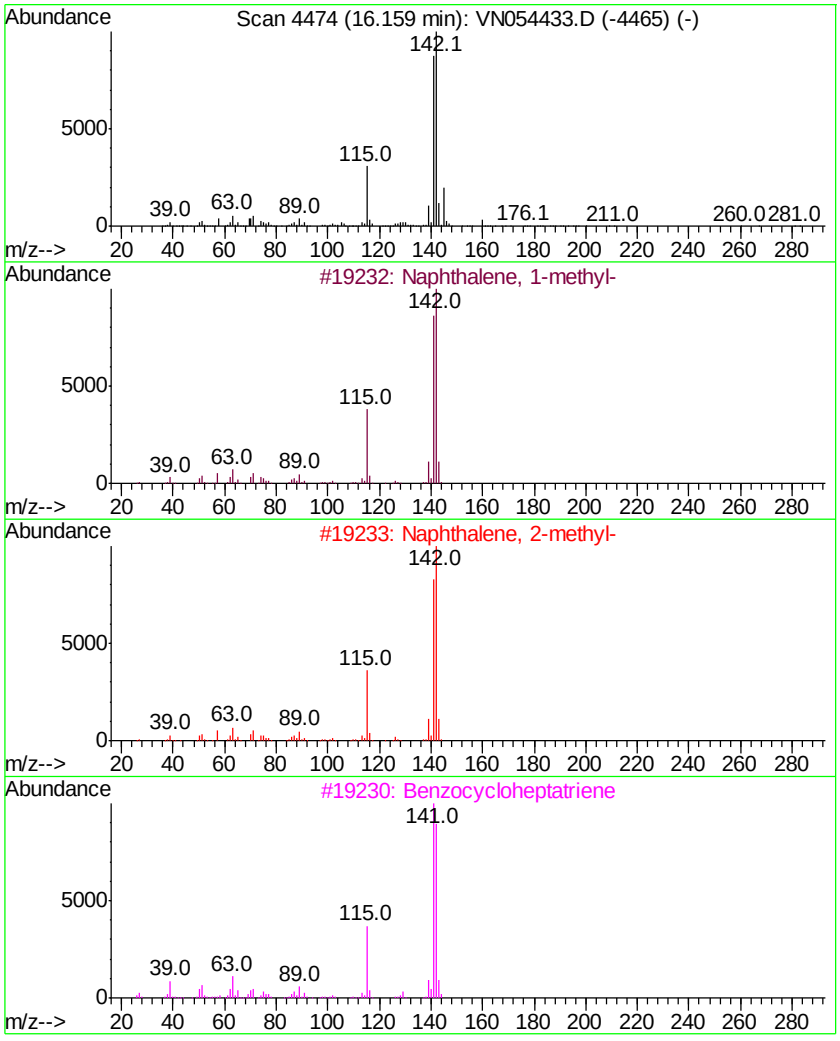
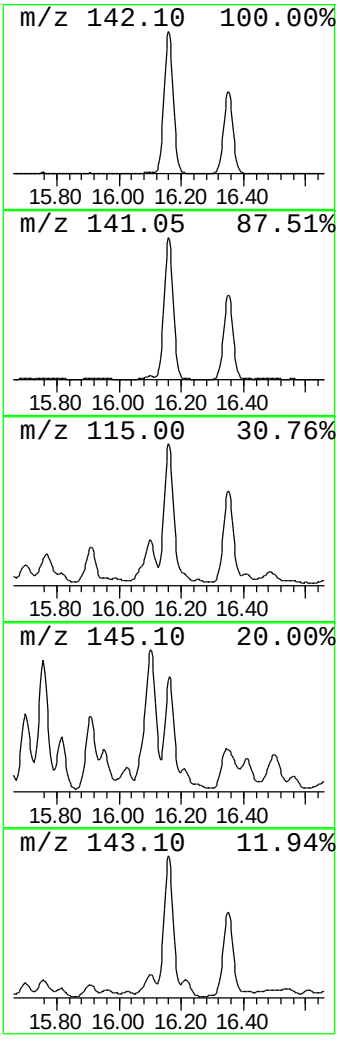
Peak Number 9 Naphthalene, 1-methyl- Concentration Rank 1

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.16	540.60 ug/l	63575000	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
2	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	90
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5	1H-Indene, 1-ethylidene-	142	C11H10	002471-83-2	53





Library Search Compound Report

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Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260

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

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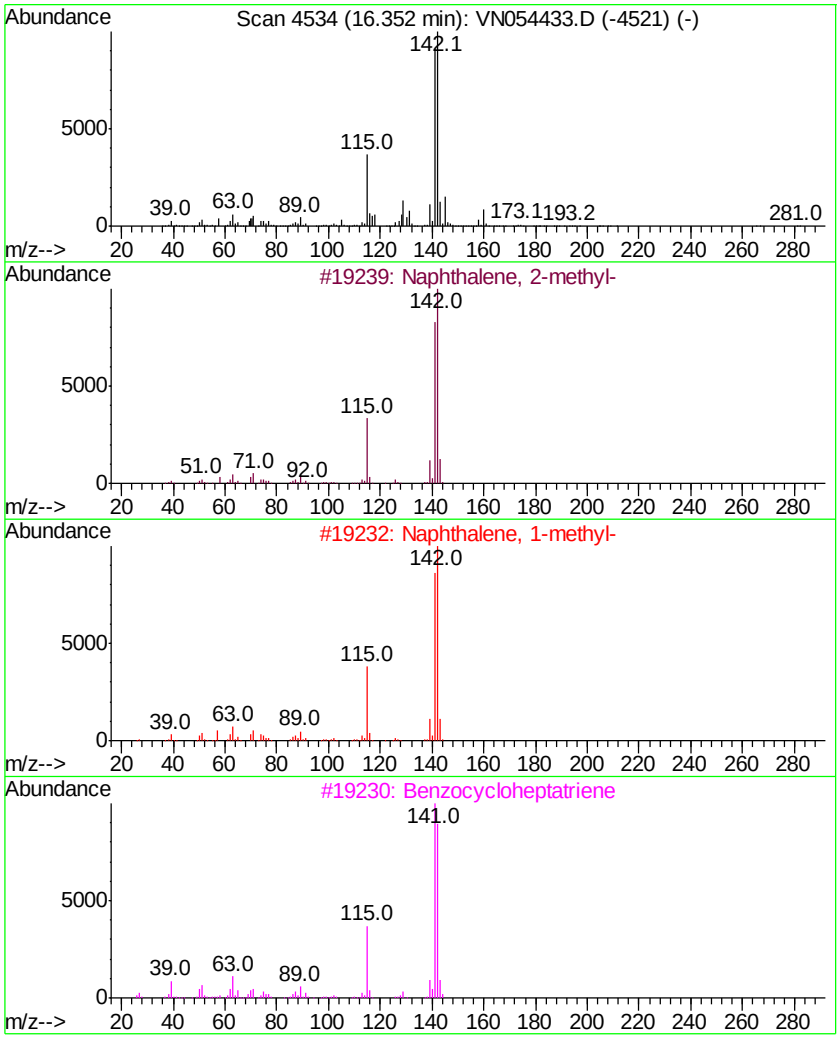
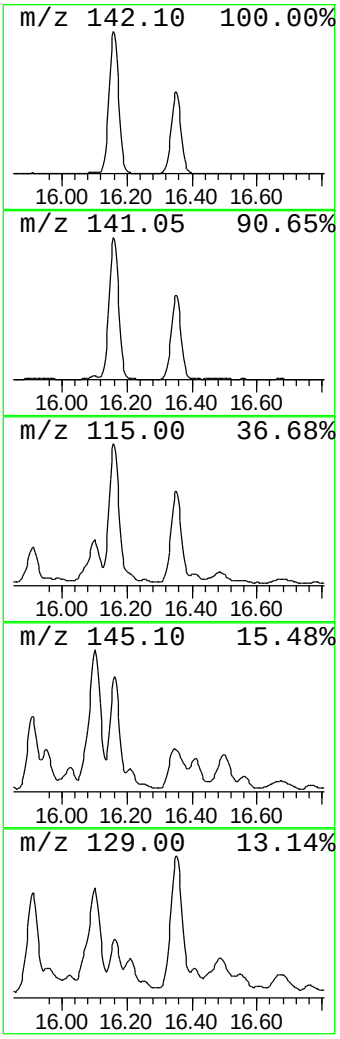
Peak Number 10 Naphthalene, 2-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
16.35	352.66 ug/l	41473000	1,4-Dichlorobenzene-d4	13.34

Hit# of 5	Tentative ID	MW	MolForm	CAS#	Qual
1	Naphthalene, 2-methyl-	142	C11H10	000091-57-6	96
2	Naphthalene, 1-methyl-	142	C11H10	000090-12-0	96
3	Benzocycloheptatriene	142	C11H10	000264-09-5	87
4	1,4-Methanonaphthalene, 1,4-dihy...	142	C11H10	004453-90-1	87
5	Spiro(tricyclo[6.2.1.0(2,7)]unde...	186	C13H14O	1000210-02-3	78





Tentatively Identified Compound (LSC) summary

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031919\
Data File : VN054433.D
Acq On : 19 Mar 2019 11:47
Operator : JC/SP
Sample : K1970-05
Misc : 5.04g/10mL/100uL/5.00mL/MSVOA_N/MEOH
ALS Vial : 5 Sample Multiplier: 1

Instrument : MSVOA_N
ClientSampleId : P-5-E.WALL-SOUTH

Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.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
Benzene, 4-ethyl-...	13.89	232.7	ug/l	27359900	4	13.34	5880090	50.0
Benzene, 1-methyl...	13.99	209.8	ug/l	24676300	4	13.34	5880090	50.0
1-Phenyl-1-butene	14.59	508.6	ug/l	59813400	4	13.34	5880090	50.0
6-Methyl-4-indanol	14.85	237.0	ug/l	27868900	4	13.34	5880090	50.0
Bicyclo[4.2.0]oct...	14.96	295.0	ug/l	34689100	4	13.34	5880090	50.0
Benzene, (1-methy...	15.61	322.6	ug/l	37933600	4	13.34	5880090	50.0
Naphthalene, 1,2,...	15.91	225.8	ug/l	26558600	4	13.34	5880090	50.0
Naphthalene, 1,2,...	16.10	277.8	ug/l	32674800	4	13.34	5880090	50.0
Naphthalene, 1-me...	16.16	540.6	ug/l	63575000	4	13.34	5880090	50.0
Naphthalene, 2-me...	16.35	352.7	ug/l	41473000	4	13.34	5880090	50.0