

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN080418\
 Data File : VN050367.D
 Acq On : 4 Aug 2018 15:21
 Operator : MD\SY
 Sample : J4328-07
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :

Integration Parameters: RTEINT.P

Integrator: RTE
 Smoothing : ON
 Sampling : 1
 Start Thrs: 0.2
 Stop Thrs : 0

Filtering: 5
 Min Area: 3 % of largest Peak
 Max Peaks: 100
 Peak Location: TOP

If leading or trailing edge < 100 prefer < Baseline drop else tangent >
 Peak separation: 5

Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N072418W.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	1.654	3	20	55	rBV	11827121	19225907	100.00%	44.918%
2	7.593	1848	1867	1878	rBV	530384	1309905	6.81%	3.060%
3	7.667	1878	1890	1912	rVB	788035	1891281	9.84%	4.419%
4	8.030	1985	2003	2031	rBV	557279	1357695	7.06%	3.172%
5	8.590	2160	2177	2203	rBV	1146592	2374472	12.35%	5.548%
6	10.091	2629	2644	2674	rBV	2066755	3807107	19.80%	8.895%
7	11.410	3036	3054	3077	rBV2	1607085	3640049	18.93%	8.504%
8	12.403	3347	3363	3378	rBV	1411255	2364565	12.30%	5.524%
9	13.345	3643	3656	3673	rBV	1484657	2503944	13.02%	5.850%
10	13.545	3700	3718	3735	rBV	624014	1171267	6.09%	2.736%
11	13.670	3746	3757	3769	rVB4	102267	200963	1.05%	0.470%
12	13.889	3816	3825	3835	rVB	183576	288186	1.50%	0.673%
13	13.998	3850	3859	3869	rVB2	316740	527599	2.74%	1.233%
14	14.210	3919	3925	3940	rVB2	169573	278130	1.45%	0.650%
15	14.586	4032	4042	4055	rBV3	239246	466329	2.43%	1.090%
16	14.959	4150	4158	4174	rVB5	157709	339849	1.77%	0.794%
17	15.191	4222	4230	4239	rBV	127501	215339	1.12%	0.503%
18	15.265	4243	4253	4280	rVB2	181958	497148	2.59%	1.162%
19	15.917	4444	4456	4471	rBV3	166229	342160	1.78%	0.799%

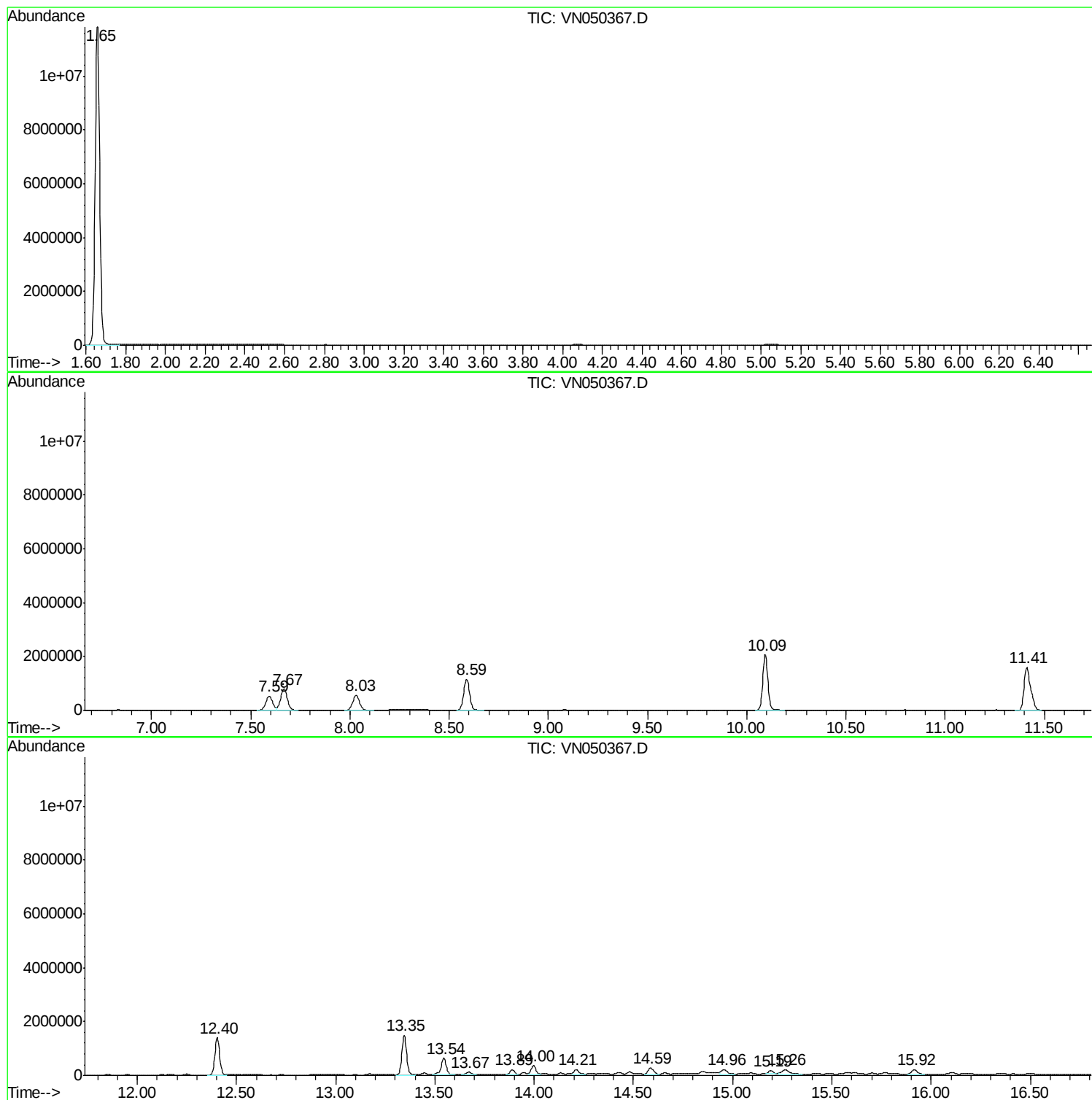
Sum of corrected areas: 42801895

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Quant Title : SW846 8260

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



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 Peak Number 1 Indane Concentration Rank 2

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.54	23.39 ug/l	1171270	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Indane	118 C9H10	000496-11-7 94
2		Benzene, cyclopropyl-	118 C9H10	000873-49-4 68
3		Tetracyclo[3.3.1.0(2,8).0(4,6)]-...	118 C9H10	1000191-13-7 64
4		Benzene, 1-propenyl-	118 C9H10	000637-50-3 62
5		trans-Cinnamyl bromide	196 C9H9Br	026146-77-0 59

 Peak Number 2 Benzene, 1-ethyl-2,3-dimethyl- Concentration Rank 8

R.T.	EstConc	Area	Relative to ISTD	R.T.
13.89	5.75 ug/l	288186	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-2,3-dimethyl-	134 C10H14	000933-98-2 96
2		o-Cymene	134 C10H14	000527-84-4 95
3		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 95
4		Benzene, 2-ethyl-1,3-dimethyl-	134 C10H14	002870-04-4 94
5		Benzene, 1-methyl-3-(1-methyleth...	134 C10H14	000535-77-3 94

 Peak Number 3 Indan, 1-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.00	10.54 ug/l	527599	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Indan, 1-methyl-	132 C10H12	000767-58-8 70
2		Benzene, 1-ethenyl-3-ethyl-	132 C10H12	007525-62-4 68
3		Benzene, 1-methyl-4-(2-propenyl)-	132 C10H12	003333-13-9 64
4		Benzene, 1-methyl-2-(2-propenyl)-	132 C10H12	001587-04-8 59
5		1H-Indene, 2,3-dihydro-4-methyl-	132 C10H12	000824-22-6 53

 Peak Number 4 Benzene, 1,2,4,5-tetramethyl- Concentration Rank 9

R.T.	EstConc	Area	Relative to ISTD	R.T.
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14.21	5.55 ug/l	278130	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,4,5-tetramethyl-	134 C10H14	000095-93-2 91
2		Benzene, 1,2,3,5-tetramethyl-	134 C10H14	000527-53-7 91
3		1,3-Cyclopentadiene, 1,2,3,4-tet...	134 C10H14	076089-59-3 91
4		Benzene, 1,2,3,4-tetramethyl-	134 C10H14	000488-23-3 91
5		Benzene, 2-ethyl-1,4-dimethyl-	134 C10H14	001758-88-9 91

 Peak Number 5 1H-Indene, 2,3-dihydro-1,3-... Concentration Rank 7

R.T.	EstConc	Area	Relative to ISTD	R.T.
14.96	6.79 ug/l	339849	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 64
2		1H-Indene, 2,3-dihydro-1,1-dimet...	146 C11H14	004912-92-9 62
3		1H-Indene, 2,3-dihydro-4,7-dimet...	146 C11H14	006682-71-9 58
4		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 58
5		Methanethioate(dimethylamino), (...	235 C13H17NOS	1000197-43-0 47

 Peak Number 6 Benzene, (3-methyl-2-butenyl)- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.26	9.93 ug/l	497148	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, (3-methyl-2-butenyl)-	146 C11H14	004489-84-3 94
2		Naphthalene, 1,2,3,4-tetrahydro-...	146 C11H14	001559-81-5 91
3		1H-Indene, 2,3-dihydro-1,6-dimet...	146 C11H14	017059-48-2 90
4		2,2-Dimethylindene, 2,3-dihydro-	146 C11H14	020836-11-7 90
5		1H-Indene, 2,3-dihydro-1,3-dimet...	146 C11H14	004175-53-5 87

 Peak Number 7 Naphthalene, 1,2,3,4-tetra... Concentration Rank 6

R.T.	EstConc	Area	Relative to ISTD	R.T.
15.92	6.83 ug/l	342160	1,4-Dichlorobenzene-d4	13.35
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual

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1	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	002809-64-5	81
2	Naphthalene, 1,2,3,4-tetrahydro-...	146	C11H14	001680-51-9	81
3	2-Propyn-1-ol, 3-(4-methylphenyl)-	146	C10H10O	016017-24-6	74
4	1H-1,3,2-Benzodiazaborole, 2-eth...	146	C8H11BN2	074663-81-3	64
5	Benzene, (1,2-dimethyl-1-propenyl)-	146	C11H14	000769-57-3	53

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Indane	13.54	23.4	ug/l	1171270	4	13.35	2503940	50.0
Benzene, 1-ethyl-...	13.89	5.8	ug/l	288186	4	13.35	2503940	50.0
Indan, 1-methyl-	14.00	10.5	ug/l	527599	4	13.35	2503940	50.0
Benzene, 1,2,4,5-...	14.21	5.5	ug/l	278130	4	13.35	2503940	50.0
1H-Indene, 2,3-di...	14.96	6.8	ug/l	339849	4	13.35	2503940	50.0
Benzene, (3-methy...	15.26	9.9	ug/l	497148	4	13.35	2503940	50.0
Naphthalene, 1,2,...	15.92	6.8	ug/l	342160	4	13.35	2503940	50.0