

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU010519\
 Data File : VU029037.D
 Acq On : 04 Jan 2019 19:06
 Operator : MD/SY
 Sample : K1061-01 10X
 Misc : 5.0mL/MSVOA U/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 TRE-COMP-1

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_U\METHOD\82U010319W.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	1.283	27	34	46	rBV	136694	133194	21.84%	2.839%
2	1.386	60	66	75	rBV2	13589	16995	2.79%	0.362%
3	1.469	84	92	108	rBV	396058	440995	72.31%	9.400%
4	2.013	252	261	271	rVB	16373	22513	3.69%	0.480%
5	2.244	320	333	346	rVB	14190	21837	3.58%	0.465%
6	2.321	348	357	373	rVB2	5131	8713	1.43%	0.186%
7	2.833	506	516	531	rBV2	3208	6404	1.05%	0.137%
8	3.530	719	733	751	rBV4	10319	25407	4.17%	0.542%
9	4.881	1137	1153	1170	rBV2	91931	208757	34.23%	4.450%
10	4.980	1170	1184	1207	rVB	140611	306232	50.21%	6.527%
11	5.305	1272	1285	1301	rBV	99738	208569	34.20%	4.446%
12	5.389	1302	1311	1327	rVB3	11238	23638	3.88%	0.504%
13	5.884	1450	1465	1480	rBV	220853	431715	70.79%	9.202%
14	5.961	1483	1489	1506	rVB4	6648	13788	2.26%	0.294%
15	7.562	1971	1987	2001	rBV	352533	609845	100.00%	12.999%
16	7.636	2001	2010	2025	rVB	38848	69514	11.40%	1.482%
17	9.086	2448	2461	2484	rBV	306858	505226	82.84%	10.769%
18	9.250	2503	2512	2523	rBV	9649	14631	2.40%	0.312%
19	9.372	2538	2550	2572	rBV	48247	81355	13.34%	1.734%
20	9.778	2665	2676	2692	rBV2	29138	45509	7.46%	0.970%
21	10.305	2829	2840	2867	rBV	227163	357723	58.66%	7.625%
22	10.588	2916	2928	2941	rBV	9208	14135	2.32%	0.301%
23	10.678	2945	2956	2975	rBV	45346	97841	16.04%	2.085%
24	10.771	2975	2985	2995	rBV	24423	35571	5.83%	0.758%
25	10.970	3037	3047	3064	rBV	22539	33836	5.55%	0.721%
26	11.147	3091	3102	3116	rBV	88896	136263	22.34%	2.904%
27	11.482	3194	3206	3224	rBV	272373	427134	70.04%	9.104%
28	11.569	3224	3233	3249	rVB	38004	57496	9.43%	1.226%
29	11.765	3279	3294	3302	rBV4	22060	49522	8.12%	1.056%
30	11.806	3302	3307	3316	rVV	10398	15884	2.60%	0.339%
31	11.874	3316	3328	3342	rVB3	17454	33023	5.41%	0.704%
32	12.054	3378	3384	3396	rVB2	5385	7640	1.25%	0.163%
33	12.144	3402	3412	3417	rBV2	11075	16015	2.63%	0.341%
34	12.173	3417	3421	3435	rVB2	10058	14331	2.35%	0.305%

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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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Title : SW846 8260

35	12.250	3435	3445	3454	rBV2	18488	27914	4.58%	0.595%
36	12.353	3468	3477	3486	rBV3	6176	10027	1.64%	0.214%
37	12.549	3531	3538	3548	rVB	5536	7227	1.19%	0.154%
38	12.649	3561	3569	3577	rBV2	10857	15576	2.55%	0.332%
39	12.700	3577	3585	3593	rBV	19172	27377	4.49%	0.584%
40	12.964	3658	3667	3680	rVB2	10319	15249	2.50%	0.325%
41	13.122	3706	3716	3733	rVB3	25302	41421	6.79%	0.883%
42	13.285	3759	3767	3777	rVB4	7935	10956	1.80%	0.234%
43	13.745	3900	3910	3928	rBV	17352	30723	5.04%	0.655%
44	14.350	4085	4098	4109	rBV5	7357	13843	2.27%	0.295%

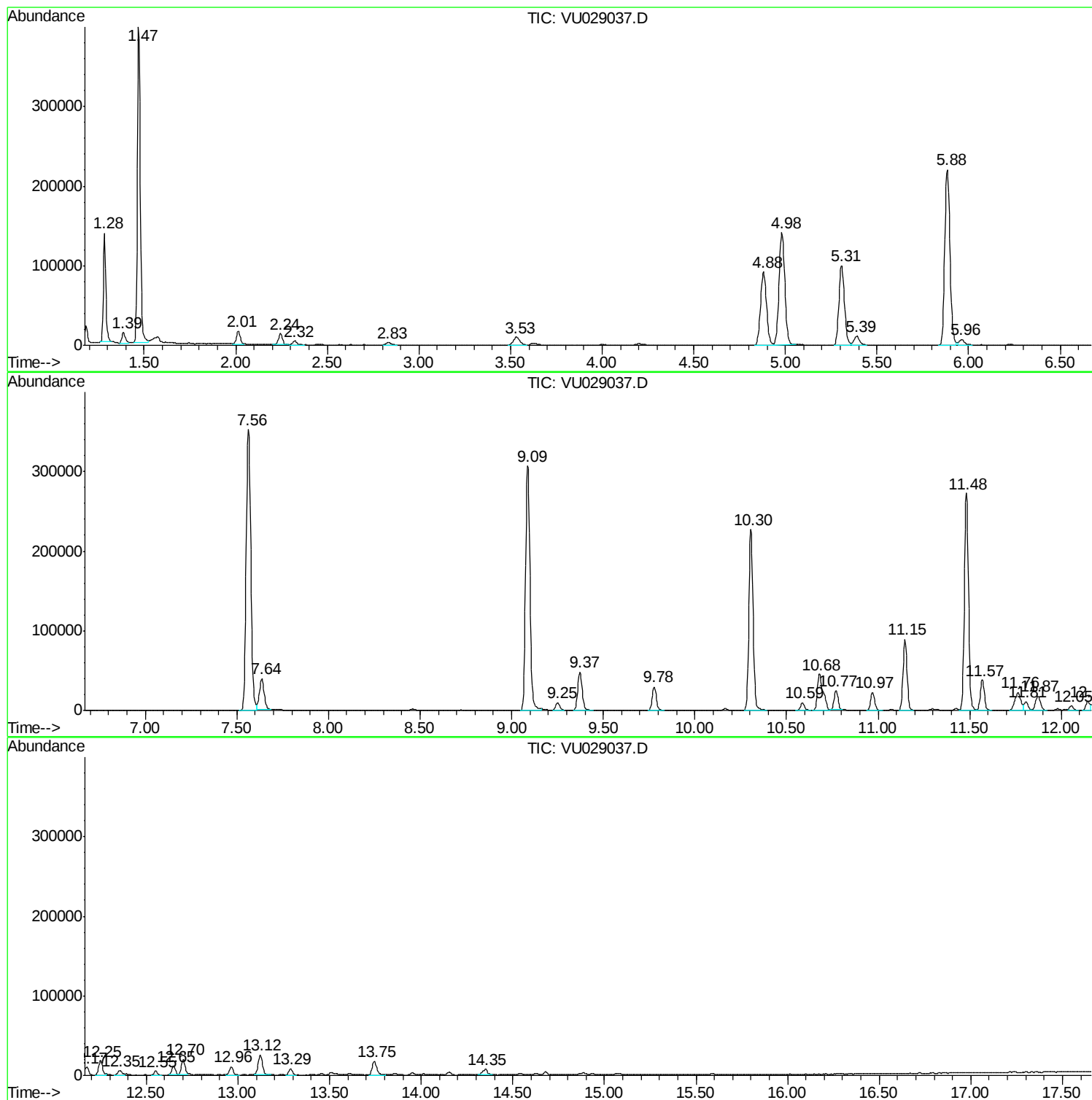
Sum of corrected areas: 4691564

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Quant Method : Z:\V0ASRV\HPCHEM1\MSVOA_U\METHOD\82U010319W.M
Quant Title : SW846 8260

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



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 Quant Title : SW846 8260

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

 Peak Number 1 Benzene, 1-ethyl-3-methyl- Concentration Rank 3

R.T.	EstConc	Area	Relative to ISTD	R.T.
10.68	11.45 ug/l	97841	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 94
2		Benzene, 1-ethyl-2-methyl-	120 C9H12	000611-14-3 91
3		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 91

 Peak Number 2 Benzene, 1,2,3-trimethyl- Concentration Rank 4

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.57	6.73 ug/l	57496	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Benzene, 1,2,3-trimethyl-	120 C9H12	000526-73-8 95
2		Benzene, 1,2,4-trimethyl-	120 C9H12	000095-63-6 94
3		Benzene, 1-ethyl-3-methyl-	120 C9H12	000620-14-4 94
4		Mesitylene	120 C9H12	000108-67-8 91
5		Benzene, 1-ethyl-4-methyl-	120 C9H12	000622-96-8 91

 Peak Number 3 Indane Concentration Rank 5

R.T.	EstConc	Area	Relative to ISTD	R.T.
11.76	5.80 ug/l	49522	1,4-Dichlorobenzene-d4	11.48
Hit# of	5	Tentative ID	MW MolForm	CAS# Qual
1		Indane	118 C9H10	000496-11-7 87
2		Benzene, 2-propenyl-	118 C9H10	000300-57-2 74
3		Benzene, cyclopropyl-	118 C9H10	000873-49-4 72
4		Benzene, 1-ethenyl-2-methyl-	118 C9H10	000611-15-4 72
5		Benzene, 1-propenyl-	118 C9H10	000637-50-3 64

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TIC Top Hit name	RT	EstConc	Units	Response	--Internal Standard--			
					#	RT	Resp	Conc
Benzene, 1-ethyl-...	10.68	11.4	ug/l	97841	4	11.48	427134	50.0
Benzene, 1,2,3-tr...	11.57	6.7	ug/l	57496	4	11.48	427134	50.0
Indane	11.76	5.8	ug/l	49522	4	11.48	427134	50.0