

Data Path : W:\HPCHEM1\MSVOA X\DATA\VX033018\
 Data File : VX000564.D
 Acq On : 30 Mar 2018 17:06
 Operator : JC/MD
 Sample : J2139-11
 Misc : 5.00g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PAD-2

Manual Integrations
 APPROVED

sam
 4/4/2018 10:29:25 AM

Quant Time: Apr 04 11:31:11 2018
 Quant Method : W:\HPCHEM1\MSVOA_X\METHOD\82X032018W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 21 04:52:42 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.67	168	150443	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.87	114	265857	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.12	117	271091	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.09	152	167226	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	6.08	65	101182	51.52	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.04%	
35) Dibromofluoromethane	5.52	113	77802	46.45	ug/l	0.00
Spiked Amount	50.000		Recovery	=	92.90%	
50) Toluene-d8	8.73	98	333039	48.00	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.00%	
62) 4-Bromofluorobenzene	11.15	95	131896	45.61	ug/l	0.00
Spiked Amount	50.000		Recovery	=	91.22%	
Target Compounds						
						Qvalue
3) Chloromethane	1.32	50	1109	0.63	ug/l	91
5) Bromomethane	1.64	94	6377	2.93	ug/l	99
20) Methylene Chloride	2.85	84	2704	1.43	ug/l	97
40) Benzene	6.15	78	126537	17.50	ug/l	100
52) Toluene	8.79	92	157155	31.08	ug/l	100
67) Ethyl Benzene	10.26	91	46392	4.54	ug/l	97
68) m/p-Xylenes	10.37	106	166820	41.45	ug/l	94
69) o-Xylene	10.71	106	61645	15.82	ug/l	98
70) Styrene	10.73	104	102741	15.94	ug/l	96
73) Isopropylbenzene	11.02	105	6018	0.54	ug/l	93
74) N-amyl acetate	6.52	43	1627	Below Cal	#	59
78) n-propylbenzene	11.37	91	5848	0.44	ug/l	98
80) 1,3,5-Trimethylbenzene	11.52	105	90096	9.67	ug/l	98
84) 1,2,4-Trimethylbenzene	11.82	105	204252	20.91	ug/l	98

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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