

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN111820\
 Data File : VN064713.D
 Acq On : 18 Nov 2020 17:48
 Operator : JC/MD
 Sample : L4706-01ME 10X
 Misc : 1.06g/10mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 COMP-30706-30707-40854ME

Manual Integrations
 APPROVED

MMDadoda
 11/19/2020 2:10:56 PM

Quant Time: Nov 18 18:11:48 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N111820W.M
 Quant Title : SW846 8260
 QLast Update : Wed Nov 18 15:29:54 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.62	168	212477	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	419620	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	437297	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	200134	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.98	65	191403	51.21	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.42%	
35) Dibromofluoromethane	7.55	113	122786	51.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	102.82%	
50) Toluene-d8	10.06	98	536622	49.44	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.88%	
62) 4-Bromofluorobenzene	12.37	95	224712	55.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	110.78%	
Target Compounds						
					Qvalue	
16) Acetone	3.77	43	5576m	11.105	ug/l	
39) Methylcyclohexane	9.04	83	13491	3.527	ug/l #	87
52) Toluene	10.12	92	1402208	189.271	ug/l	99
67) Ethyl Benzene	11.48	91	86202	5.256	ug/l	99
68) m/p-Xylenes	11.59	106	162832	27.518	ug/l	100
69) o-Xylene	11.92	106	58943	10.301	ug/l	99
73) Isopropylbenzene	12.22	105	20869	1.469	ug/l	97
78) n-propylbenzene	12.56	91	91071	5.503	ug/l	99
80) 1,3,5-Trimethylbenzene	12.70	105	98042	8.306	ug/l	100
84) 1,2,4-Trimethylbenzene	13.01	105	245930	20.414	ug/l	99
85) sec-Butylbenzene	13.14	105	6195m	0.509	ug/l	
86) p-Isopropyltoluene	13.26	119	4324	0.386	ug/l	87
89) n-Butylbenzene	13.59	91	7396	0.708	ug/l #	81

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

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