

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN072018\
 Data File : VN049971.D
 Acq On : 20 Jul 2018 18:57
 Operator : MD\SY
 Sample : PB111289ZHE#11
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PB111289ZHE#11

Manual Integrations
 APPROVED

MMDadoda
 7/23/2018 2:04:11 PM

Quant Time: Jul 21 02:23:32 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N071818W.M
 Quant Title : SW846 8260
 QLast Update : Thu Jul 19 01:58:39 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	554266	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	949775	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	803416	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	288394	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	413211	54.24	ug/l	0.00
Spiked Amount 50.000			Recovery	=	108.48%	
35) Dibromofluoromethane	7.59	113	358774	45.00	ug/l	0.00
Spiked Amount 50.000			Recovery	=	90.00%	
50) Toluene-d8	10.09	98	1263597	43.50	ug/l	0.00
Spiked Amount 50.000			Recovery	=	87.00%	
62) 4-Bromofluorobenzene	12.40	95	344315	34.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	68.82%	
Target Compounds						
16) Acetone	3.82	43	1854890	940.43	ug/l	Qvalue 100
18) Methyl Acetate	4.33	43	31530	4.33	ug/l	99
20) Methylene Chloride	4.55	84	1870953	249.43	ug/l	99
25) 2-Butanone	6.84	43	25879m	9.18	ug/l	
29) Tetrahydrofuran	7.23	42	15169	8.54	ug/l	94
43) Isopropyl Acetate	8.17	43	346169	27.32	ug/l #	91
59) 2-Hexanone	10.77	43	20838	4.38	ug/l #	57

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

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