

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN082318\
 Data File : VN050856.D
 Acq On : 23 Aug 2018 16:36
 Operator : MD\SY
 Sample : J4617-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 N48969

Manual Integrations
 APPROVED

MMDadoda
 8/24/2018 10:40:58 AM

Quant Time: Aug 24 11:58:08 2018
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N082118W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 22 05:38:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	672973	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	997715	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	846817	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	300953	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	405026	49.99	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.98%	
35) Dibromofluoromethane	7.59	113	386628	48.71	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.42%	
50) Toluene-d8	10.09	98	1360577	45.85	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.70%	
62) 4-Bromofluorobenzene	12.40	95	376679	38.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	76.40%	
Target Compounds						
16) Acetone	3.82	43	385139	199.32	ug/l	98
29) Tetrahydrofuran	7.22	42	747303	399.22	ug/l	99
64) Tetrachloroethene	10.63	164	3092	0.38	ug/l	89
67) Ethyl Benzene	11.51	91	10388	0.29	ug/l	99
68) m/p-Xylenes	11.62	106	4349	0.32	ug/l	88
69) o-Xylene	11.95	106	8324	0.63	ug/l	98

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

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