

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN090619\
 Data File : VN057734.D
 Acq On : 6 Sep 2019 23:22
 Operator : JC/SP
 Sample : K4580-33
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PAR-12-C2-(3.5)

Manual Integrations
 APPROVED

MMDadoda
 9/10/2019 9:56:58 AM

Quant Time: Sep 07 00:16:55 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N090619W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 06 03:46:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.65	168	191656	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.58	114	326025	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	293800	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	151121	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.01	65	254850	54.66	ug/l	0.00
Spiked Amount 50.000			Recovery	=	109.32%	
35) Dibromofluoromethane	7.58	113	147789	49.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.54%	
50) Toluene-d8	10.08	98	573445	52.20	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.40%	
62) 4-Bromofluorobenzene	12.40	95	208215	45.63	ug/l	0.00
Spiked Amount 50.000			Recovery	=	91.26%	
Target Compounds						
16) Acetone	3.81	43	9978	6.954	ug/l	92
20) Methylene Chloride	4.52	84	5480m	1.766	ug/l	
43) Isopropyl Acetate	8.15	43	71627	8.666	ug/l	98

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

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