

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN122418\
 Data File : VN053189.D
 Acq On : 24 Dec 2018 17:00
 Operator : MD\SY
 Sample : J6511-14
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 19 Sample Multiplier: 28

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-59(0-5)-GRAB

Manual Integrations
 APPROVED

MMDadoda
 12/26/2018 11:20:59 AM

Quant Time: Dec 25 00:32:23 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N121818W.M
 Quant Title : SW846 8260
 QLast Update : Tue Dec 18 13:03:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1210940	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	1884030	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	1632071	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	622356	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	606836	49.38	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.76%	
35) Dibromofluoromethane	7.59	113	521080	60.55	ug/l	0.00
Spiked Amount 50.000			Recovery	=	121.10%	
50) Toluene-d8	10.09	98	2270694	49.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.30%	
62) 4-Bromofluorobenzene	12.40	95	719921	44.39	ug/l	0.00
Spiked Amount 50.000			Recovery	=	88.78%	
Target Compounds						
16) Acetone	3.82	43	26774	11.112	ug/l	96
18) Methyl Acetate	4.33	43	5743m	1.507	ug/l	
20) Methylene Chloride	4.55	84	32453	2.469	ug/l	97
25) 2-Butanone	6.84	43	6062	1.499	ug/l	98
43) Isopropyl Acetate	8.17	43	52257	1.665	ug/l #	82

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

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