

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN061820\
 Data File : VN061940.D
 Acq On : 19 Jun 2020 8:30
 Operator : JC/MD
 Sample : L2813-12
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 56 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EO-01-061720

Manual Integrations
 APPROVED

MMDadoda
 6/19/2020 1:02:06 PM

Quant Time: Jun 19 12:02:43 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N061620W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 17 01:19:44 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	142738	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	272503	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	268768	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	104562	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	100860	52.69	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.38%	
35) Dibromofluoromethane	7.55	113	86769	49.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	99.54%	
50) Toluene-d8	10.06	98	343957	50.19	ug/l	0.00
Spiked Amount 50.000			Recovery	=	100.38%	
62) 4-Bromofluorobenzene	12.37	95	113812	48.41	ug/l	0.00
Spiked Amount 50.000			Recovery	=	96.82%	
Target Compounds						
16) Acetone	3.78	43	10403	18.397	ug/l	96
18) Methyl Acetate	4.28	43	3660	2.248	ug/l #	85
20) Methylene Chloride	4.51	84	6476	3.515	ug/l	91
25) 2-Butanone	6.80	43	3348m	4.058	ug/l	
43) Isopropyl Acetate	8.13	43	29441	9.001	ug/l #	90
51) 4-Methyl-2-Pentanone	9.95	43	3114	1.611	ug/l	87

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

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