

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN100720\
 Data File : VN063940.D
 Acq On : 7 Oct 2020 14:08
 Operator : JC/MD
 Sample : L4279-02
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 S-1

Manual Integrations
 APPROVED

MMDadoda
 10/8/2020 10:08:43 AM

Quant Time: Oct 08 16:19:20 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N093020W.M
 Quant Title : SW846 8260
 QLast Update : Thu Oct 01 01:46:38 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.63	168	215770	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.55	114	387169	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.38	117	403069	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.32	152	156792	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.99	65	187181	52.51	ug/l	0.00
Spiked Amount 50.000			Recovery	=	105.02%	
35) Dibromofluoromethane	7.55	113	136396	50.77	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.54%	
50) Toluene-d8	10.06	98	498268	48.93	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.86%	
62) 4-Bromofluorobenzene	12.38	95	187207	48.90	ug/l	0.00
Spiked Amount 50.000			Recovery	=	97.80%	
Target Compounds						
16) Acetone	3.78	43	118928	108.380	ug/l	97
17) Carbon Disulfide	4.01	76	96730	16.941	ug/l	100
20) Methylene Chloride	4.51	84	5054	2.105	ug/l #	76
25) 2-Butanone	6.79	43	16122m	10.363	ug/l	
43) Isopropyl Acetate	8.13	43	9446	1.470	ug/l #	90

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

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