

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042220\
 Data File : VN061145.D
 Acq On : 22 Apr 2020 22:10
 Operator : JC/MD
 Sample : L2359-04
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-1B-S

Quant Time: Apr 23 08:43:03 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042220W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 22 15:26:40 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.40	168	193644	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.35	114	323783	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.21	117	314646	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.16	152	153035	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	7.77	65	122891	47.86	ug/l	0.00
Spiked Amount 50.000			Recovery	=	95.72%	
35) Dibromofluoromethane	7.31	113	90827	49.15	ug/l	0.00
Spiked Amount 50.000			Recovery	=	98.30%	
50) Toluene-d8	9.88	98	402079	50.78	ug/l	0.00
Spiked Amount 50.000			Recovery	=	101.56%	
62) 4-Bromofluorobenzene	12.21	95	155140	52.32	ug/l	0.00
Spiked Amount 50.000			Recovery	=	104.64%	
Target Compounds						
16) Acetone	3.55	43	20094	9.577	ug/l	98
20) Methylene Chloride	4.24	84	3917	1.746	ug/l	94
40) Benzene	7.78	78	7429	0.793	ug/l	98
43) Isopropyl Acetate	7.92	43	54431	10.105	ug/l	99

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

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