

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042419\
 Data File : VN055229.D
 Acq On : 24 Apr 2019 15:44
 Operator : JC/SP
 Sample : K2200-08
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EO-02-042319-A

Quant Time: Apr 25 06:48:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N042219W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 23 05:52:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	386246	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	660331	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	603353	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	183619	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	263096	50.33	ug/l	0.00
Spiked Amount	50.000		Recovery	=	100.66%	
35) Dibromofluoromethane	7.59	113	213507	47.59	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.18%	
50) Toluene-d8	10.09	98	818376	49.24	ug/l	0.00
Spiked Amount	50.000		Recovery	=	98.48%	
62) 4-Bromofluorobenzene	12.40	95	256132	48.66	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.32%	
Target Compounds						
16) Acetone	3.82	43	17028	9.853	ug/l	92
17) Carbon Disulfide	4.05	76	4268	0.384	ug/l #	92
18) Methyl Acetate	4.34	43	8898	1.298	ug/l	94
20) Methylene Chloride	4.55	84	126764	25.514	ug/l	93
30) Chloroform	7.37	83	4606	0.575	ug/l	88
43) Isopropyl Acetate	8.17	43	22869	2.855	ug/l #	78

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

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