

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN042419\
 Data File : VN055223.D
 Acq On : 24 Apr 2019 13:11
 Operator : JC/SP
 Sample : K2542-16
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TP-11

Quant Time: Apr 25 05:13:24 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	364235	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	638056	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	579657	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	180857	50.00	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.03	65	254306	51.58	ug/l	0.00
Spiked Amount	50.000		Recovery	=	103.16%	
35) Dibromofluoromethane	7.59	113	206431	47.62	ug/l	0.00
Spiked Amount	50.000		Recovery	=	95.24%	
50) Toluene-d8	10.09	98	780269	48.58	ug/l	0.00
Spiked Amount	50.000		Recovery	=	97.16%	
62) 4-Bromofluorobenzene	12.40	95	244956	48.16	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.32%	
Target Compounds						
8) Diethyl Ether	3.41	74	3985	1.782	ug/l	87
16) Acetone	3.82	43	10321	6.333	ug/l	91
17) Carbon Disulfide	4.05	76	5081	0.485	ug/l	100
20) Methylene Chloride	4.55	84	70956	15.144	ug/l	99
30) Chloroform	7.37	83	6723	0.890	ug/l	88
43) Isopropyl Acetate	8.17	43	16945	2.189	ug/l #	88

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

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