

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN071422\
 Data File : VN073463.D
 Acq On : 14 Jul 2022 16:14
 Operator : JC\MD
 Sample : N3715-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 GWOW-BR-03-280-300-071322

Quant Time: Jul 15 02:02:12 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N062822W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jun 29 01:54:03 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.016	168	266090	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	456075	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	406441	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	185380	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	177597	46.234	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	92.460%
35) Dibromofluoromethane	7.951	113	146242	48.196	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	96.400%
50) Toluene-d8	10.381	98	471552	43.756	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	87.520%
62) 4-Bromofluorobenzene	12.674	95	196878	47.538	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	95.080%
Target Compounds						
					Qvalue	
16) Acetone	4.228	43	13141	6.095	ug/l #	79
25) 2-Butanone	7.263	43	3923	1.143	ug/l #	82
40) Benzene	8.392	78	13143	0.912	ug/l	94
51) 4-Methyl-2-Pentanone	10.269	43	17561	2.731	ug/l #	86
52) Toluene	10.439	92	21281	2.390	ug/l	94

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

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