

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062722\
 Data File : VN073170.D
 Acq On : 27 Jun 2022 18:29
 Operator : JC\MD
 Sample : N3436-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 31677

Quant Time: Jun 28 06:39:01 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	334712	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	530212	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	485535	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	247404	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.375	65	198545	43.875	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.760%
35) Dibromofluoromethane	7.951	113	170374	51.988	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.980%
50) Toluene-d8	10.380	98	564504	46.292	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.580%
62) 4-Bromofluorobenzene	12.668	95	249043	56.795	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	113.580%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	108748	43.718	ug/l	93
17) Carbon Disulfide	4.457	76	150059	18.276	ug/l	99
25) 2-Butanone	7.269	43	32178	8.152	ug/l #	89
29) Tetrahydrofuran	7.622	42	8273	3.174	ug/l	95
40) Benzene	8.392	78	57460	4.123	ug/l	96
52) Toluene	10.439	92	37151	4.342	ug/l	99
68) m/p-Xylenes	11.892	106	14360	2.193	ug/l	92

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

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