

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN080322\
 Data File : VN073710.D
 Acq On : 03 Aug 2022 14:35
 Operator : JC\MD
 Sample : N3971-15
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-23R

Quant Time: Aug 04 05:34:11 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N072822W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jul 29 02:20:25 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	360927	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	568506	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	546617	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.616	152	261481	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	210544	47.867	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	95.740%
35) Dibromofluoromethane	7.951	113	193580	54.226	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	108.460%
50) Toluene-d8	10.380	98	651228	46.358	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.720%
62) 4-Bromofluorobenzene	12.674	95	247053	52.195	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.380%
Target Compounds						Qvalue
16) Acetone	4.234	43	40059	16.956	ug/l	95
19) Methyl tert-butyl Ether	5.534	73	21832	1.866	ug/l #	84
31) Cyclohexane	8.016	56	51189	7.817	ug/l #	91
37) Ethyl Acetate	7.339	43	52733	7.870	ug/l	98
39) Methylcyclohexane	9.398	83	2392	0.393	ug/l	97
40) Benzene	8.392	78	58920	3.566	ug/l	100
52) Toluene	10.445	92	6100	0.604	ug/l	88
67) Ethyl Benzene	11.786	91	56954	2.908	ug/l	96
68) m/p-Xylenes	11.904	106	8823	1.181	ug/l	99
69) o-Xylene	12.221	106	1556	0.214	ug/l	89
73) Isopropylbenzene	12.521	105	33651	1.758	ug/l	99
78) n-propylbenzene	12.863	91	33337	1.548	ug/l	93
84) 1,2,4-Trimethylbenzene	13.310	105	4435	0.286	ug/l	94
95) Naphthalene	15.439	128	12865	0.866	ug/l	99

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

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