

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042122\
 Data File : VX028245.D
 Acq On : 21 Apr 2022 16:46
 Operator : JC/MD
 Sample : N2459-12
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 31582

Quant Time: Apr 22 06:15:31 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X041922W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 19 13:38:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	381000	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	655831	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	632039	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	299908	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	277954	55.951	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	111.900%
35) Dibromofluoromethane	5.385	113	254159	55.260	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	110.520%
50) Toluene-d8	8.653	98	921419	52.161	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	104.320%
62) 4-Bromofluorobenzene	11.079	95	336147	49.768	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.540%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	6968	3.214	ug/l	94
25) 2-Butanone	4.574	43	1084	0.368	ug/l	97
52) Toluene	8.720	92	5529	0.453	ug/l	95
67) Ethyl Benzene	10.195	91	7497	0.332	ug/l	99
68) m/p-Xylenes	10.299	106	11096	1.207	ug/l	99
69) o-Xylene	10.646	106	5981	0.652	ug/l	99
84) 1,2,4-Trimethylbenzene	11.756	105	7222	0.415	ug/l	98

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

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