

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX052422\
 Data File : VX028919.D
 Acq On : 24 May 2022 14:22
 Operator : JC/MD
 Sample : N2958-01
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TT162S1-20220516

Quant Time: May 25 01:48:54 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051222W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 12 14:58:53 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	346830	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	614751	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	581312	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	272677	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	233825	49.182	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	98.360%
35) Dibromofluoromethane	5.391	113	201340	49.988	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	99.980%
50) Toluene-d8	8.653	98	724991	47.743	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.480%
62) 4-Bromofluorobenzene	11.079	95	282870	49.873	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	99.740%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	3339	1.636	ug/l	94
24) 1,1-Dichloroethane	3.611	63	7694	1.076	ug/l	95
44) Trichloroethene	7.129	130	52089	11.047	ug/l	98
64) Tetrachloroethene	9.281	164	3677	0.838	ug/l #	87

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

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