

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122022\
 Data File : VX033503.D
 Acq On : 20 Dec 2022 13:22
 Operator : JC/MD
 Sample : N6149-02 100X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PARAM-1

Quant Time: Dec 21 04:14:03 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 07 12:43:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	118810	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	187785	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	166661	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	79041	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	35928	47.137	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	94.280%
35) Dibromofluoromethane	5.391	113	41442	49.066	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.140%
50) Toluene-d8	8.647	98	81520	46.137	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	92.280%
62) 4-Bromofluorobenzene	11.079	95	63327	46.874	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.740%
Target Compounds						
					Qvalue	
16) Acetone	2.392	43	6145	10.895	ug/l	100
52) Toluene	8.720	92	6054	1.950	ug/l	98
67) Ethyl Benzene	10.195	91	15430	2.624	ug/l	97
68) m/p-Xylenes	10.299	106	33340	14.574	ug/l	100
69) o-Xylene	10.646	106	17884	7.909	ug/l	99

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

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