

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122022\
 Data File : VX033511.D
 Acq On : 20 Dec 2022 16:26
 Operator : JC/MD
 Sample : N6149-09 100X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PARS-1

Quant Time: Dec 21 04:15:39 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	120616	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	194212	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	175026	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	82976	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	36892	47.677	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	95.360%
35) Dibromofluoromethane	5.385	113	43903	50.259	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	100.520%
50) Toluene-d8	8.647	98	88106	48.215	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.420%
62) 4-Bromofluorobenzene	11.079	95	69758	49.926	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	99.860%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	7775	32.865	ug/l	99
16) Acetone	2.392	43	6916	12.079	ug/l	98
52) Toluene	8.720	92	11981	3.732	ug/l	95
67) Ethyl Benzene	10.195	91	14979	2.426	ug/l	98
68) m/p-Xylenes	10.299	106	26652	11.094	ug/l	97
69) o-Xylene	10.640	106	14588	6.143	ug/l	98

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

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