

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122022\
 Data File : VX033516.D
 Acq On : 20 Dec 2022 18:21
 Operator : JC/MD
 Sample : N6149-09 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PARS-1

Quant Time: Dec 21 04:16:38 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	96383	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	153306	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	196663	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	95134	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	27410	44.329	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	88.660%
35) Dibromofluoromethane	5.385	113	33803	49.023	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.040%
50) Toluene-d8	8.647	98	96924	67.193	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	134.380%#
62) 4-Bromofluorobenzene	11.079	95	78022	70.740	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	141.480%#
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.965	59	61379	324.682	ug/l #	76
16) Acetone	2.380	43	47962	104.824	ug/l	100
18) Methyl Acetate	2.703	43	10455	7.893	ug/l	100
19) Methyl tert-butyl Ether	3.111	73	14624	4.832	ug/l	96
20) Methylene Chloride	2.788	84	1690	1.656	ug/l	92
25) 2-Butanone	4.556	43	6489	9.880	ug/l #	86
40) Benzene	6.044	78	7987	2.038	ug/l #	90
43) Isopropyl Acetate	6.348	43	3910	1.912	ug/l #	92
52) Toluene	8.720	92	122266	48.244	ug/l	100
67) Ethyl Benzene	10.195	91	168396	24.271	ug/l	98
68) m/p-Xylenes	10.299	106	295921	109.626	ug/l	96
69) o-Xylene	10.640	106	160843	60.282	ug/l	96
84) 1,2,4-Trimethylbenzene	11.750	105	17528	3.059	ug/l	98

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

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