

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122022\
 Data File : VX033518.D
 Acq On : 20 Dec 2022 19:07
 Operator : JC/MD
 Sample : N6149-07
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FAIR-1

Quant Time: Dec 21 04:17:04 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	146159	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	193563	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	190642	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.073	152	59634	50.000	ug/l	0.05
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	40033	42.695	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	85.380%
35) Dibromofluoromethane	5.391	113	50169	57.625	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	115.260%
50) Toluene-d8	8.653	98	101398	55.674	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	111.340%
62) 4-Bromofluorobenzene	11.091	95	82012	58.893	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	117.780%
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.983	59	87044	303.635	ug/l	98
16) Acetone	2.380	43	394877	569.116	ug/l	97
18) Methyl Acetate	2.703	43	32771	16.315	ug/l	98
20) Methylene Chloride	2.794	84	11970	7.736	ug/l	86
25) 2-Butanone	4.556	43	251721	252.743	ug/l	94
49) 1,4-Dioxane	7.702	88	25371	835.032	ug/l #	80
52) Toluene	8.720	92	14524	4.539	ug/l	99
67) Ethyl Benzene	10.201	91	8002	1.190	ug/l	97
68) m/p-Xylenes	10.305	106	14241	5.442	ug/l	98
69) o-Xylene	10.646	106	8480	3.279	ug/l	91
80) 1,3,5-Trimethylbenzene	11.463	105	5388	1.498	ug/l	98

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

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