

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090321\
 Data File : VX024110.D
 Acq On : 03 Sep 2021 18:14
 Operator : JC/MD
 Sample : M3646-16
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 93-94

Quant Time: Sep 04 01:32:15 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X081721W.M
 Quant Title : SW846 8260
 QLast Update : Tue Aug 17 11:57:51 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	133004	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	221974	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	208645	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	91649	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.970	65	102564	55.251	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	= 110.500%		
35) Dibromofluoromethane	5.397	113	80640	54.887	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 109.780%		
50) Toluene-d8	8.653	98	285218	51.006	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	= 102.020%		
62) 4-Bromofluorobenzene	11.085	95	104269	50.653	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	= 101.300%		
Target Compounds						
					Qvalue	
3) Chloromethane	1.294	50	882	0.573	ug/l #	87
5) Bromomethane	1.618	94	1334	1.412	ug/l	91
11) Tert butyl alcohol	2.971	59	4198	11.183	ug/l	100
16) Acetone	2.386	43	11465	13.813	ug/l	90
20) Methylene Chloride	2.800	84	1447	0.913	ug/l	91
25) 2-Butanone	4.587	43	2219	1.765	ug/l	91
40) Benzene	6.050	78	1673	0.275	ug/l #	90

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

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