

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042825\
 Data File : VX045958.D
 Acq On : 28 Apr 2025 15:49
 Operator : JC/MD
 Sample : Q1891-08
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MH-D

Quant Time: Apr 29 01:34:57 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 02 03:11:43 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	64553	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	127577	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	120116	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	52395	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	64528	54.661	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 109.320%	
35) Dibromofluoromethane	5.385	113	45785	50.582	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 101.160%	
50) Toluene-d8	8.647	98	161060	50.979	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 101.960%	
62) 4-Bromofluorobenzene	11.079	95	60576	52.638	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 105.280%	
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	54796	113.040	ug/l	97
18) Methyl Acetate	2.709	43	2078	1.862	ug/l	95
20) Methylene Chloride	2.788	84	3000	3.306	ug/l	95
25) 2-Butanone	4.587	43	6499	9.159	ug/l	97
27) cis-1,2-Dichloroethene	4.489	96	5760	5.972	ug/l	90
43) Isopropyl Acetate	6.348	43	9899	4.241	ug/l	98

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

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