

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX051523\
 Data File : VX035657.D
 Acq On : 15 May 2023 12:07
 Operator : JC/MD
 Sample : 02643-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FSND-PRO-07-99-05032023

Quant Time: May 16 04:56:49 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X051123W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 11 13:57:54 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 05/16/2023
 Supervised By : Mahesh Dadoda 05/16/2023

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	205332	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	379328	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	345949	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	162934	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	172557	48.172	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	96.340%		
35) Dibromofluoromethane	5.385	113	123438	48.252	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.500%		
50) Toluene-d8	8.647	98	456430	48.297	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	96.600%		
62) 4-Bromofluorobenzene	11.079	95	193375	50.819	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	101.640%		
Target Compounds					Qvalue	
12) 1,1-Dichloroethene	2.319	96	845	0.370	ug/l	95
16) Acetone	2.404	43	21277	13.820	ug/l #	87
17) Carbon Disulfide	2.514	76	23363	4.233	ug/l	99
18) Methyl Acetate	2.715	43	1842m	0.324	ug/l	
21) trans-1,2-Dichloroethene	3.093	96	633	0.253	ug/l #	72
25) 2-Butanone	4.568	43	16240	7.620	ug/l	98
30) Chloroform	5.105	83	1767	0.336	ug/l	91
44) Trichloroethene	7.129	130	1785750	623.449	ug/l	93
59) 2-Hexanone	9.427	43	296460	93.038	ug/l	98

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

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