

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX071522\
 Data File : VX030086.D
 Acq On : 15 Jul 2022 13:06
 Operator : JC/MD
 Sample : N3735-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 GWOW-BR-03-370-390-071422

Quant Time: Jul 16 06:01:37 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X071222W.M
 Quant Title : SW846 8260
 QLast Update : Wed Jul 13 09:05:23 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	163044	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	289745	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	239930	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	93821	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	97052	48.904	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	97.800%		
35) Dibromofluoromethane	5.391	113	86946	46.973	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.940%		
50) Toluene-d8	8.653	98	325673	46.782	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	93.560%		
62) 4-Bromofluorobenzene	11.085	95	102179	42.295	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	84.600%		
Target Compounds						
16) Acetone	2.386	43	6158	6.585	ug/l	89
25) 2-Butanone	4.568	43	2851	1.790	ug/l #	68
40) Benzene	6.050	78	8312	1.105	ug/l	97
51) 4-Methyl-2-Pentanone	8.580	43	3307	1.082	ug/l	94
52) Toluene	8.726	92	8997	1.883	ug/l	96
64) Tetrachloroethene	9.275	164	1002	0.520	ug/l #	76

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

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