

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072122\
 Data File : VX030170.D
 Acq On : 21 Jul 2022 18:26
 Operator : JC/MD
 Sample : N3765-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB-04-07132022

Quant Time: Jul 22 04:59:35 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	5.556	168	167716	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	302959	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	271306	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	111944	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	113839	55.765	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 111.520%			
35) Dibromofluoromethane	5.391	113	93070	48.088	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 96.180%			
50) Toluene-d8	8.653	98	358355	49.231	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 98.460%			
62) 4-Bromofluorobenzene	11.085	95	119260	47.213	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 94.420%			
Target Compounds					Qvalue	
16) Acetone	2.385	43	12021	12.497	ug/l	94
17) Carbon Disulfide	2.513	76	33775	6.738	ug/l	96
25) 2-Butanone	4.574	43	4136	2.524	ug/l	99
44) Trichloroethene	7.122	130	610	0.265	ug/l	42
52) Toluene	8.726	92	1238	0.248	ug/l	86
64) Tetrachloroethene	9.281	164	3162	1.452	ug/l	91

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

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