

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072922\
 Data File : VX030306.D
 Acq On : 29 Jul 2022 16:32
 Operator : JC/MD
 Sample : N3861-09
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-17

Quant Time: Aug 01 05:15:18 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X072522W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jul 25 16:49:17 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	138383	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	244235	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	249632	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	138346	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	153766	55.283	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	110.560%
35) Dibromofluoromethane	5.391	113	121200	51.911	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.820%
50) Toluene-d8	8.653	98	457482	50.597	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	101.200%
62) 4-Bromofluorobenzene	11.085	95	161658	48.175	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	96.360%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	9303	6.533	ug/l	95
31) Cyclohexane	5.483	56	20972	5.193	ug/l #	86
39) Methylcyclohexane	7.391	83	7912	1.824	ug/l	92
40) Benzene	6.050	78	24984	2.363	ug/l	97
78) n-propylbenzene	11.311	91	3124	0.242	ug/l	100

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

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