

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX072122\
 Data File : VX030168.D
 Acq On : 21 Jul 2022 17:39
 Operator : JC/MD
 Sample : N3781-02DL 250X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ORA-0019DL

Quant Time: Jul 22 06:59:30 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	168630	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	304377	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.061	117	267736	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	113646	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.964	65	112025	54.579	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 109.160%			
35) Dibromofluoromethane	5.391	113	91309	46.958	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 93.920%			
50) Toluene-d8	8.653	98	357407	48.872	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 97.740%			
62) 4-Bromofluorobenzene	11.085	95	119035	46.904	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 93.800%			
Target Compounds					Qvalue	
13) Acrolein	2.252	56	735	6.967	ug/l	92
16) Acetone	2.386	43	117188	121.166	ug/l	98
25) 2-Butanone	4.580	43	2204	1.338	ug/l	99
51) 4-Methyl-2-Pentanone	8.580	43	2811	0.875	ug/l	98
52) Toluene	8.720	92	22992	4.580	ug/l	94
68) m/p-Xylenes	10.305	106	3113	0.878	ug/l	99
69) o-Xylene	10.640	106	1227	0.346	ug/l	97

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

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