

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091522\
 Data File : VX031367.D
 Acq On : 15 Sep 2022 20:29
 Operator : JC/MD
 Sample : N4508-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 BR-01-180-200-090122

Quant Time: Sep 16 04:58:01 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 01:50:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	92249	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	153406	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	139842	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	64030	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	100023	44.174	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	88.340%		
35) Dibromofluoromethane	5.385	113	82893	50.269	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	100.540%		
50) Toluene-d8	8.653	98	309246	49.994	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	99.980%		
62) 4-Bromofluorobenzene	11.079	95	99634	43.986	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	87.980%		
Target Compounds					Qvalue	
12) 1,1-Dichloroethene	2.319	96	13719	10.652	ug/l	94
16) Acetone	2.386	43	2760	2.855	ug/l #	86
17) Carbon Disulfide	2.514	76	83381	27.458	ug/l	100
24) 1,1-Dichloroethane	3.617	63	2074	0.664	ug/l #	86
27) cis-1,2-Dichloroethene	4.495	96	4341	2.288	ug/l	81
30) Chloroform	5.105	83	17202	5.199	ug/l	89
44) Trichloroethene	7.129	130	7063	4.274	ug/l	88
52) Toluene	8.720	92	6476	1.580	ug/l	95

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

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