

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX070523\
 Data File : VX036417.D
 Acq On : 05 Jul 2023 20:28
 Operator : JC/MD
 Sample : 03507-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 24 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CP-1

Quant Time: Jul 06 04:57:59 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X062323W.M
 Quant Title : SW846 8260
 QLast Update : Fri Jun 23 15:03:28 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	170208	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	289403	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	258005	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	117484	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	109433	44.853	ug/l	0.00
Spiked Amount	50.000	Range 78 - 117	Recovery	=	89.700%	
35) Dibromofluoromethane	5.391	113	91992	48.208	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	=	96.420%	
50) Toluene-d8	8.647	98	347230	48.718	ug/l	0.00
Spiked Amount	50.000	Range 92 - 112	Recovery	=	97.440%	
62) 4-Bromofluorobenzene	11.079	95	123901	49.354	ug/l	0.00
Spiked Amount	50.000	Range 83 - 123	Recovery	=	98.700%	
Target Compounds						
						Qvalue
16) Acetone	2.404	43	225034	215.788	ug/l	98
18) Methyl Acetate	2.715	43	4516	1.053	ug/l	97
20) Methylene Chloride	2.794	84	9422	4.196	ug/l	90
25) 2-Butanone	4.574	43	119485	79.744	ug/l	96
37) Ethyl Acetate	4.721	43	51720	17.840	ug/l	97
43) Isopropyl Acetate	6.342	43	194257	42.003	ug/l	97
51) 4-Methyl-2-Pentanone	8.574	43	531767	178.465	ug/l #	87

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

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