

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU060722\
 Data File : VU048907.D
 Acq On : 07 Jun 2022 14:15
 Operator : SY/MD
 Sample : N3165-09
 Misc : 5.0mL/MSVOA_U/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 60BR-20220531

Quant Time: Jun 08 04:51:57 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\82U052622W.M
 Quant Title : SW846 8260
 QLast Update : Fri May 27 06:42:31 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.372	168	252242	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.250	114	403189	50.000	ug/l	0.00
63) Chlorobenzene-d5	9.417	117	368204	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.812	152	159784	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.700	65	186129	49.134	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	98.260%
35) Dibromofluoromethane	5.289	113	140733	48.926	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.860%
50) Toluene-d8	7.896	98	502806	48.764	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.520%
62) 4-Bromofluorobenzene	10.632	95	171217	43.907	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	87.820%
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.385	85	2017	0.698	ug/l	99
16) Acetone	2.629	43	7819	4.050	ug/l	91
21) trans-1,2-Dichloroethene	3.350	96	1450	0.494	ug/l #	78
27) cis-1,2-Dichloroethene	4.665	96	184572	54.299	ug/l	88
37) Ethyl Acetate	4.800	43	509423	95.987	ug/l	99
44) Trichloroethene	6.543	130	839265	262.837	ug/l	100
60) Dibromochloromethane	8.806	129	1903	0.519	ug/l	100
71) Bromoform	10.295	173	964	0.353	ug/l #	94

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

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