

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN010322\
 Data File : VN070447.D
 Acq On : 3 Jan 2022 19:24
 Operator : JC/MD
 Sample : M5338-09
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TW-2

Quant Time: Jan 04 01:37:20 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N122721W.M
 Quant Title : SW846 8260
 QLast Update : Mon Dec 27 18:47:12 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.086	168	964400	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.965	114	1631349	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.744	117	1422672	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.675	152	479019	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.437	65	696213	50.209	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	= 100.420%	
35) Dibromofluoromethane	8.021	113	480107	48.641	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	= 97.280%	
50) Toluene-d8	10.440	98	1904959	47.244	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	= 94.480%	
62) 4-Bromofluorobenzene	12.731	95	671175	46.086	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	= 92.180%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	4.189	96	29734	3.027	ug/l	92
16) Acetone	4.293	43	25651	2.850	ug/l	99
24) 1,1-Dichloroethane	6.391	63	99266	4.414	ug/l	99
25) 2-Butanone	7.348	43	22050	1.946	ug/l #	89
27) cis-1,2-Dichloroethene	7.335	96	4149	0.334	ug/l #	66
30) Chloroform	7.820	83	39304	1.812	ug/l	91
32) 1,1,1-Trichloroethane	8.013	97	25046	1.369	ug/l #	50
38) Carbon Tetrachloride	8.206	117	10220	0.655	ug/l	92
52) Toluene	10.505	92	10494	0.346	ug/l	95
64) Tetrachloroethene	10.977	164	30631	2.973	ug/l	96
67) Ethyl Benzene	11.843	91	17783	0.322	ug/l	97
68) m/p-Xylenes	11.950	106	21766	1.081	ug/l	98
84) 1,2,4-Trimethylbenzene	13.366	105	16188	0.472	ug/l	95

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

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