

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN050622\
 Data File : VN072357.D
 Acq On : 06 May 2022 16:11
 Operator : JC\MD
 Sample : N2725-08 20X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 12 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TRE-1281

Quant Time: May 09 03:13:54 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N042622W.M
 Quant Title : SW846 8260
 QLast Update : Tue Apr 26 18:50:11 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.081	168	242544	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	465523	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	537232	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.669	152	215858	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	188004	50.490	ug/l	0.00
Spiked Amount	50.000	Range 61 - 141	Recovery	= 100.980%		
35) Dibromofluoromethane	8.022	113	143445	48.142	ug/l	0.00
Spiked Amount	50.000	Range 69 - 133	Recovery	= 96.280%		
50) Toluene-d8	10.439	98	607434	51.882	ug/l	0.00
Spiked Amount	50.000	Range 65 - 126	Recovery	= 103.760%		
62) 4-Bromofluorobenzene	12.727	95	254552	65.503	ug/l	0.00
Spiked Amount	50.000	Range 58 - 135	Recovery	= 131.000%		
Target Compounds						
						Qvalue
16) Acetone	4.281	43	518509	322.980	ug/l	94
20) Methylene Chloride	5.087	84	1634	0.531	ug/l #	59
25) 2-Butanone	7.340	43	45124	17.649	ug/l	90
37) Ethyl Acetate	7.416	43	25673	4.457	ug/l	97
51) 4-Methyl-2-Pentanone	10.328	43	8886	1.635	ug/l #	85
52) Toluene	10.498	92	12461	1.549	ug/l	99
68) m/p-Xylenes	11.939	106	1953	0.278	ug/l	85

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

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