

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN050622\
 Data File : VN072355.D
 Acq On : 06 May 2022 15:23
 Operator : JC\MD
 Sample : N2725-06 20X
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 TRE-1279

Quant Time: May 09 03:13:23 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	8.086	168	242891	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	450781	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	486071	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.674	152	200636	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	180884	48.508	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	97.020%
35) Dibromofluoromethane	8.016	113	140983	48.863	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	97.720%
50) Toluene-d8	10.439	98	582589	51.387	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	102.780%
62) 4-Bromofluorobenzene	12.727	95	226003	60.058	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	120.120%
Target Compounds						
					Qvalue	
16) Acetone	4.287	43	124392	77.373	ug/l	92
20) Methylene Chloride	5.104	84	3265	1.059	ug/l #	82
25) 2-Butanone	7.334	43	29063	11.351	ug/l #	89
37) Ethyl Acetate	7.410	43	196772	35.281	ug/l	97
51) 4-Methyl-2-Pentanone	10.327	43	15002	2.851	ug/l #	80
52) Toluene	10.498	92	16922	2.173	ug/l	99
68) m/p-Xylenes	11.945	106	2173	0.342	ug/l	89

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

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