

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN042922\
 Data File : VN072274.D
 Acq On : 29 Apr 2022 16:58
 Operator : JC\MD
 Sample : N2601-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 BP-MW150S1-GW-245-247

Quant Time: Apr 30 03:15:31 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	185897	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	8.963	114	319907	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	274140	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	81335	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.434	65	167439	58.670	ug/l	0.00
Spiked Amount	50.000	Range	61 - 141	Recovery	=	117.340%
35) Dibromofluoromethane	8.016	113	109253	53.357	ug/l	0.00
Spiked Amount	50.000	Range	69 - 133	Recovery	=	106.720%
50) Toluene-d8	10.439	98	385613	47.927	ug/l	0.00
Spiked Amount	50.000	Range	65 - 126	Recovery	=	95.860%
62) 4-Bromofluorobenzene	12.727	95	117166	43.873	ug/l	0.00
Spiked Amount	50.000	Range	58 - 135	Recovery	=	87.740%
Target Compounds						
						Qvalue
16) Acetone	4.293	43	4478	3.639	ug/l	92
31) Cyclohexane	8.081	56	4835	1.268	ug/l #	39
44) Trichloroethene	9.210	130	56802	26.414	ug/l	86
64) Tetrachloroethene	10.969	164	2555	1.548	ug/l #	84

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

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