

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061822\
 Data File : VN072934.D
 Acq On : 18 Jun 2022 12:54
 Operator : JC\MD
 Sample : N3329-02RE
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PALMYRA-RT-73-COMP-1RE

Quant Time: Jun 20 08:03:35 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N061722W.M
 Quant Title : SW846 8260
 QLast Update : Sat Jun 18 06:22:45 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	290465	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	516411	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	443335	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	188121	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	174946	44.549	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	89.100%
35) Dibromofluoromethane	7.951	113	138457	43.378	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	86.760%
50) Toluene-d8	10.374	98	455478	38.349	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	76.700%#
62) 4-Bromofluorobenzene	12.663	95	160529	37.587	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	75.180%#
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	50376	23.337	ug/l	89
20) Methylene Chloride	5.022	84	76913	22.840	ug/l	96
25) 2-Butanone	7.269	43	22935	6.695	ug/l #	88
43) Isopropyl Acetate	8.498	43	42871	4.341	ug/l #	88

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

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