

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN061822\
 Data File : VN072935.D
 Acq On : 18 Jun 2022 13:19
 Operator : JC\MD
 Sample : N3329-04RE
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 PALMYRA-RT-73-COMP-2RE

Quant Time: Jun 20 08:03:45 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	292179	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	526152	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	448608	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	187644	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	175966	44.546	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	89.100%
35) Dibromofluoromethane	7.951	113	133583	41.076	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	82.160%
50) Toluene-d8	10.374	98	421737	34.851	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	69.700%#
62) 4-Bromofluorobenzene	12.668	95	147379	33.869	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	67.740%#
Target Compounds						
					Qvalue	
16) Acetone	4.234	43	49658	22.869	ug/l #	88
18) Methyl Acetate	4.793	43	9683	1.675	ug/l #	87
20) Methylene Chloride	5.016	84	74095	21.864	ug/l	93
25) 2-Butanone	7.269	43	22957	6.663	ug/l #	89
43) Isopropyl Acetate	8.492	43	41837	4.158	ug/l #	88

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

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