

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN053124\
 Data File : VN082415.D
 Acq On : 31 May 2024 17:59
 Operator : JC\MD
 Sample : P2662-04
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WCS-TPR

Quant Time: May 31 23:01:17 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N051524W.M
 Quant Title : SW846 8260
 QLast Update : Thu May 16 05:21:06 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.229	168	149549	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	278189	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	280416	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	118868	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.582	65	131952	55.878	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.760%
35) Dibromofluoromethane	8.171	113	94120	54.999	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	110.000%
50) Toluene-d8	10.570	98	370683	56.498	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	113.000%
62) 4-Bromofluorobenzene	12.853	95	124502	53.166	ug/l	0.00
Spiked Amount	50.000	Range	64 - 133	Recovery	=	106.340%
Target Compounds						
						Qvalue
16) Acetone	4.430	43	81374	84.004	ug/l	94
20) Methylene Chloride	5.277	84	14702	7.522	ug/l #	65
25) 2-Butanone	7.488	43	12885	8.773	ug/l #	83
43) Isopropyl Acetate	8.694	43	252780	44.414	ug/l #	72
51) 4-Methyl-2-Pentanone	10.447	43	18094	5.595	ug/l #	72

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

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