

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN081624\
 Data File : VN083362.D
 Acq On : 16 Aug 2024 18:29
 Operator : JC\MD
 Sample : P3629-03
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-1

Quant Time: Aug 17 02:55:46 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N080724W.M
 Quant Title : SW846 8260
 QLast Update : Thu Aug 08 06:30:41 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	127510	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.100	114	260931	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.865	117	275793	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	127330	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	109143	60.135	ug/l	0.00
Spiked Amount	50.000	Range 74 - 125	Recovery	= 120.280%		
35) Dibromofluoromethane	8.171	113	87666	53.827	ug/l	0.00
Spiked Amount	50.000	Range 75 - 124	Recovery	= 107.660%		
50) Toluene-d8	10.571	98	332318	54.699	ug/l	0.00
Spiked Amount	50.000	Range 86 - 113	Recovery	= 109.400%		
62) 4-Bromofluorobenzene	12.853	95	135656	57.275	ug/l	0.00
Spiked Amount	50.000	Range 77 - 121	Recovery	= 114.540%		
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	5.530	59	2278	6.040	ug/l #	77
16) Acetone	4.430	43	44818	63.107	ug/l	96
20) Methylene Chloride	5.283	84	7145	4.371	ug/l	94
25) 2-Butanone	7.494	43	9690	8.886	ug/l #	84
40) Benzene	8.606	78	57892	7.888	ug/l	99
43) Isopropyl Acetate	8.688	43	361771	78.066	ug/l #	73
51) 4-Methyl-2-Pentanone	10.447	43	14314	5.487	ug/l	89
52) Toluene	10.629	92	70391	15.179	ug/l	98
67) Ethyl Benzene	11.965	91	21023	1.880	ug/l	95
68) m/p-Xylenes	12.065	106	30204	7.212	ug/l	92
69) o-Xylene	12.400	106	14182	3.433	ug/l	96
84) 1,2,4-Trimethylbenzene	13.482	105	11864	1.320	ug/l	83
95) Naphthalene	15.641	128	203501	23.805	ug/l	99

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

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