

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN020524\
 Data File : VN080905.D
 Acq On : 05 Feb 2024 18:38
 Operator : JC\MD
 Sample : P1362-12
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-3

Quant Time: Feb 06 04:46:56 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N012924W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jan 29 12:54:23 2024
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By : John Carlone 02/06/2024
 Supervised By : Mahesh Dadoda 02/06/2024

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	78699	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	194145	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	217515	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	85858	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	68879	50.591	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 101.180%			
35) Dibromofluoromethane	8.165	113	52679	44.351	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 88.700%			
50) Toluene-d8	10.571	98	219068	47.122	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 94.240%			
62) 4-Bromofluorobenzene	12.853	95	88477	54.615	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 109.220%			
Target Compounds						Qvalue
3) Chloromethane	2.359	50	2059	1.172	ug/l #	70
16) Acetone	4.430	43	12164	24.635	ug/l	99
18) Methyl Acetate	5.006	43	3178	1.847	ug/l #	69
20) Methylene Chloride	5.283	84	6022m	4.646	ug/l	
25) 2-Butanone	7.483	43	5781	6.164	ug/l #	80
37) Ethyl Acetate	7.559	43	3660	1.610	ug/l #	92
43) Isopropyl Acetate	8.694	43	124511	30.092	ug/l #	89

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

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