

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN062522\
 Data File : VN073108.D
 Acq On : 25 Jun 2022 14:16
 Operator : JC\MD
 Sample : N3431-11
 Misc : 4.88g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-14B-2-3-GR

Manual Integrations
 APPROVED

Reviewed By : John Carlone 06/27/2022
 Supervised By : Mahesh Dadoda 07/08/2022

Quant Time: Jun 27 03:39:23 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	8.010	168	361391	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	565932	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	518907	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	393380	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.363	65	200912	41.121	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	82.240%		
35) Dibromofluoromethane	7.945	113	171081	48.909	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	97.820%		
50) Toluene-d8	10.374	98	715719	54.988	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	109.980%		
62) 4-Bromofluorobenzene	12.668	95	423185	90.417	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	180.840%#		
Target Compounds						
					Qvalue	
17) Carbon Disulfide	4.381	76	10159	1.146	ug/l	97
18) Methyl Acetate	4.775	43	78820	11.021	ug/l	93
39) Methylcyclohexane	9.386	83	980239	171.818	ug/l	95
73) Isopropylbenzene	12.515	105	225505	7.410	ug/l	98
85) sec-Butylbenzene	13.439	105	334081	10.939	ug/l #	76
89) n-Butylbenzene	13.880	91	151895	7.523	ug/l	98

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

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