

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
 Data File : VN074241.D
 Acq On : 01 Sep 2022 14:41
 Operator : JC\MD
 Sample : N4421-09
 Misc : 3.36g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-PALM-SP-02-GR

Quant Time: Sep 01 16:16:10 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N083022W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 31 07:02:40 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.010	168	243554	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	449617	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.680	117	588730	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	377915	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.363	65	169172	44.777	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	89.560%		
35) Dibromofluoromethane	7.945	113	135377	46.271	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	92.540%		
50) Toluene-d8	10.374	98	532585	46.223	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	92.440%		
62) 4-Bromofluorobenzene	12.668	95	320701	91.527	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	183.060%#		
Target Compounds						
					Qvalue	
18) Methyl Acetate	4.769	43	159673	23.918	ug/l #	88
39) Methylcyclohexane	9.386	83	32765	5.880	ug/l #	90
40) Benzene	8.380	78	40441	2.713	ug/l	94
52) Toluene	10.439	92	16699	1.901	ug/l	93
67) Ethyl Benzene	11.786	91	43028	1.810	ug/l	98
68) m/p-Xylenes	11.892	106	41913	4.746	ug/l	80
69) o-Xylene	12.221	106	12014	1.365	ug/l	89
78) n-propylbenzene	12.863	91	104899	2.843	ug/l	100
80) 1,3,5-Trimethylbenzene	12.998	105	130337	4.952	ug/l	70
84) 1,2,4-Trimethylbenzene	13.304	105	210515	8.287	ug/l	92
85) sec-Butylbenzene	13.439	105	85114	2.695	ug/l #	76

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

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