

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090122\
 Data File : VN074240.D
 Acq On : 01 Sep 2022 14:16
 Operator : JC\MD
 Sample : N4421-04
 Misc : 4.77g/5mL/100uL/5.00mL/MSVOA_N/MEOH
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 WC-PALM-SP-01-GR

Quant Time: Sep 01 16:15:57 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	208191	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.898	114	388589	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	409071	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.610	152	303554	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.363	65	138428	42.863	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.720%
35) Dibromofluoromethane	7.945	113	112924	44.658	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	89.320%
50) Toluene-d8	10.375	98	437734	43.958	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	87.920%
62) 4-Bromofluorobenzene	12.668	95	234990	77.598	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	155.200%#
Target Compounds						
					Qvalue	
18) Methyl Acetate	4.775	43	63184	9.937	ug/l #	87
39) Methylcyclohexane	9.392	83	31932	6.630	ug/l	96
67) Ethyl Benzene	11.786	91	20639	1.249	ug/l	98
68) m/p-Xylenes	11.886	106	22075	3.597	ug/l	100
69) o-Xylene	12.221	106	21462	3.510	ug/l	94
73) Isopropylbenzene	12.521	105	64422	2.478	ug/l	98
78) n-propylbenzene	12.857	91	117124	3.953	ug/l	93
80) 1,3,5-Trimethylbenzene	12.998	105	172870	8.177	ug/l	100
84) 1,2,4-Trimethylbenzene	13.304	105	473601	23.210	ug/l	99
85) sec-Butylbenzene	13.439	105	194426	7.664	ug/l	78
86) p-Isopropyltoluene	13.557	119	117488	5.952	ug/l	95
89) n-Butylbenzene	13.880	91	80576	4.687	ug/l #	75

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

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