

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX101922\
 Data File : VX032211.D
 Acq On : 19 Oct 2022 16:37
 Operator : JC/MD
 Sample : N5170-01MEDL 10X
 Misc : 5.07g/10mL/100uL/5.00mL/MSVOA_X/MEOH
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1639

Quant Time: Oct 19 17:10:47 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X101922W.M
 Quant Title : SW846 8260
 QLast Update : Wed Oct 19 13:54:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.549	168	144152	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	228337	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	203852	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	100743	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	90499	48.259	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	96.520%
35) Dibromofluoromethane	5.385	113	73164	46.046	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	92.100%
50) Toluene-d8	8.646	98	271771	47.516	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.040%
62) 4-Bromofluorobenzene	11.079	95	97629	46.613	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	93.220%
Target Compounds						
					Qvalue	
16) Acetone	2.379	43	149998	246.479	ug/l	100
39) Methylcyclohexane	7.385	83	7277	2.980	ug/l	96
52) Toluene	8.720	92	170594	44.455	ug/l	98
67) Ethyl Benzene	10.195	91	19174	2.591	ug/l	100
68) m/p-Xylenes	10.299	106	25791	8.826	ug/l	99
69) o-Xylene	10.640	106	12835	4.585	ug/l	97
73) Isopropylbenzene	10.963	105	9056	1.260	ug/l	97
78) n-propylbenzene	11.305	91	32267	3.920	ug/l	99
80) 1,3,5-Trimethylbenzene	11.451	105	36381	5.859	ug/l	100
84) 1,2,4-Trimethylbenzene	11.750	105	137185	22.725	ug/l	99
85) sec-Butylbenzene	11.890	105	19207	2.558	ug/l	87
89) n-Butylbenzene	12.335	91	44343	7.881	ug/l #	57

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

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