

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091522\
 Data File : VX031360.D
 Acq On : 15 Sep 2022 17:45
 Operator : JC/MD
 Sample : N4588-01ME
 Misc : 5.07g/10mL/100uL/5.0mL/MSVOA_X/MEOH
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 AUD-1647ME

Quant Time: Sep 16 04:55:48 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X091422W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 01:50:07 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	117354	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.763	114	189330	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	159934	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	72081	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	168537	58.510	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery = 117.020%			
35) Dibromofluoromethane	5.385	113	129832	63.794	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery = 127.580%#			
50) Toluene-d8	8.647	98	509299	66.713	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery = 133.420%#			
62) 4-Bromofluorobenzene	11.085	95	160216	57.311	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery = 114.620%			
Target Compounds					Qvalue	
16) Acetone	2.434	43	225037	183.002	ug/l	99
25) 2-Butanone	4.617	43	15884	7.831	ug/l	99
39) Methylcyclohexane	7.373	83	10496	3.421	ug/l	96
52) Toluene	8.720	92	852132	168.404	ug/l	99
67) Ethyl Benzene	10.195	91	9631	1.255	ug/l	100
68) m/p-Xylenes	10.299	106	9548	3.277	ug/l	99
69) o-Xylene	10.646	106	4334	1.465	ug/l	94
78) n-propylbenzene	11.305	91	7906	1.265	ug/l	98
80) 1,3,5-Trimethylbenzene	11.451	105	5339	1.197	ug/l	97
84) 1,2,4-Trimethylbenzene	11.756	105	20500	4.624	ug/l	99

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

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