

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX122923\
 Data File : VX039547.D
 Acq On : 29 Dec 2023 15:54
 Operator : JC/MD
 Sample : 06038-10
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 T-703

Quant Time: Dec 30 02:03:01 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X121323W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 14 01:28:15 2023
 Response via : Initial Calibration

Manual Integrations APPROVED

Reviewed By : Semsettin Yesilyurt 01/01/2024
 Supervised By : Mahesh Dadoda 01/01/2024

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	87810	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	180475	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	179514	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	75396	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	78895	49.045	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	98.100%		
35) Dibromofluoromethane	5.385	113	58564	46.971	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	93.940%		
50) Toluene-d8	8.647	98	232501	50.229	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.460%		
62) 4-Bromofluorobenzene	11.079	95	94443	49.451	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	98.900%		
Target Compounds						Qvalue
11) Tert butyl alcohol	2.983	59	5640	13.962	ug/l	97
16) Acetone	2.380	43	27389	6.581	ug/l	94
20) Methylene Chloride	2.788	84	6356	4.740	ug/l	98
25) 2-Butanone	4.568	43	17433	12.559	ug/l	97
37) Ethyl Acetate	4.733	43	3349m	1.378	ug/l	
40) Benzene	6.044	78	7406	1.331	ug/l	96
43) Isopropyl Acetate	6.342	43	114219	28.896	ug/l	99
67) Ethyl Benzene	10.195	91	12579	1.725	ug/l	96
95) Naphthalene	13.774	128	129901	23.558	ug/l	99

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

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