

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010324\
 Data File : VX039613.D
 Acq On : 03 Jan 2024 19:23
 Operator : JC/MD
 Sample : P1002-08DL 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 1216DL

Quant Time: Jan 04 03:38:47 2024
 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

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	91547	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	198192	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	214243	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	109750	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	79806	47.586	ug/l	0.00
Spiked Amount 50.000	Range 78 - 117		Recovery =	95.180%		
35) Dibromofluoromethane	5.385	113	64940	47.429	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	94.860%		
50) Toluene-d8	8.647	98	256287	50.418	ug/l	0.00
Spiked Amount 50.000	Range 92 - 112		Recovery =	100.840%		
62) 4-Bromofluorobenzene	11.079	95	119653	57.051	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	114.100%		
Target Compounds						Qvalue
11) Tert butyl alcohol	2.953	59	7597	18.039	ug/l	97
16) Acetone	2.374	43	70515	53.917	ug/l	92
17) Carbon Disulfide	2.514	76	2902	0.842	ug/l	96
25) 2-Butanone	4.556	43	30787	21.274	ug/l	96
40) Benzene	6.044	78	3817	0.625	ug/l	96
51) 4-Methyl-2-Pentanone	8.574	43	3396	1.228	ug/l	100
52) Toluene	8.720	92	5916	1.604	ug/l	96
59) 2-Hexanone	9.439	43	3137	1.314	ug/l	90
67) Ethyl Benzene	10.195	91	4421	0.508	ug/l	98
68) m/p-Xylenes	10.299	106	6259	2.005	ug/l	86
69) o-Xylene	10.640	106	2904	0.962	ug/l	95
84) 1,2,4-Trimethylbenzene	11.750	105	4153	0.561	ug/l	100
95) Naphthalene	13.774	128	11776	1.467	ug/l	98

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

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