

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX120822\
 Data File : VX033321.D
 Acq On : 08 Dec 2022 14:18
 Operator : JC/MD
 Sample : N5912-01DL 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 FMC-TOTE-1207DL

Quant Time: Dec 09 00:47:46 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X120722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Dec 07 12:43:54 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	154795	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	243427	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	215619	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	101059	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	40520	40.803	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	81.600%
35) Dibromofluoromethane	5.391	113	43050	39.319	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	78.640%
50) Toluene-d8	8.653	98	93575	40.854	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	81.700%#
62) 4-Bromofluorobenzene	11.079	95	71062	40.577	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	81.160%#
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.946	59	7760	25.559	ug/l #	74
16) Acetone	2.392	43	170881	232.542	ug/l	99
25) 2-Butanone	4.574	43	4993	4.734	ug/l #	87
30) Chloroform	5.092	83	3392	1.083	ug/l	88
52) Toluene	8.720	92	3230	0.803	ug/l	97
67) Ethyl Benzene	10.195	91	2449	0.322	ug/l #	89
68) m/p-Xylenes	10.305	106	3784	1.279	ug/l	95
69) o-Xylene	10.646	106	2556	0.874	ug/l	95
80) 1,3,5-Trimethylbenzene	11.451	105	2471	0.405	ug/l	100
84) 1,2,4-Trimethylbenzene	11.756	105	7276	1.195	ug/l	100

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

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