

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX062222\
 Data File : VX029690.D
 Acq On : 22 Jun 2022 11:44
 Operator : JC/MD
 Sample : N3286-02 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW-2

Quant Time: Jun 23 01:08:38 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X061822W.M
 Quant Title : SW846 8260
 QLast Update : Mon Jun 20 01:31:01 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	271111	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	455997	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	395050	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	175169	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	151091	42.101	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	84.200%
35) Dibromofluoromethane	5.391	113	135556	45.616	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	91.240%
50) Toluene-d8	8.653	98	513147	47.268	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	94.540%
62) 4-Bromofluorobenzene	11.085	95	182331	44.537	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	89.080%
Target Compounds						
					Qvalue	
17) Carbon Disulfide	2.514	76	2121	0.298	ug/l	96
29) Tetrahydrofuran	5.032	42	1918	1.273	ug/l	89
31) Cyclohexane	5.477	56	10147	2.106	ug/l	92
39) Methylcyclohexane	7.385	83	4608	0.919	ug/l	97
67) Ethyl Benzene	10.195	91	333209	23.895	ug/l	97
68) m/p-Xylenes	10.305	106	534488	98.287	ug/l	100
69) o-Xylene	10.646	106	24600	4.579	ug/l	96
73) Isopropylbenzene	10.963	105	19002	1.447	ug/l	100
78) n-propylbenzene	11.305	91	47758	3.229	ug/l	99
80) 1,3,5-Trimethylbenzene	11.451	105	78479	7.095	ug/l	100
84) 1,2,4-Trimethylbenzene	11.756	105	379206	34.374	ug/l	100
95) Naphthalene	13.780	128	163113	13.275	ug/l	99

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

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