

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX091522\
 Data File : VX031394.D
 Acq On : 16 Sep 2022 08:08
 Operator : JC/MD
 Sample : N4537-02
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 MW2

Quant Time: Sep 16 18:55:05 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	93378	50.000	ug/l	-0.01
34) 1,4-Difluorobenzene	6.763	114	153393	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	136089	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	64802	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	97768	42.656	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	85.320%
35) Dibromofluoromethane	5.385	113	80506	48.825	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	97.660%
50) Toluene-d8	8.653	98	303353	49.046	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	98.100%
62) 4-Bromofluorobenzene	11.079	95	97595	43.089	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	86.180%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	50485	51.596	ug/l	99
25) 2-Butanone	4.562	43	14174	8.782	ug/l	93
30) Chloroform	5.092	83	5664	1.691	ug/l	98
40) Benzene	6.037	78	4807	0.742	ug/l	98
67) Ethyl Benzene	10.195	91	1939	0.297	ug/l	97
73) Isopropylbenzene	10.963	105	1373	0.287	ug/l	96
78) n-propylbenzene	11.305	91	1167	0.208	ug/l	82
86) p-Isopropyltoluene	12.006	119	12156	2.969	ug/l	97

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

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