

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN090123\
 Data File : VN079130.D
 Acq On : 01 Sep 2023 18:11
 Operator : JC\MD
 Sample : 04232-09
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 SB-11

Quant Time: Sep 01 23:18:36 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N081523W.M
 Quant Title : SW846 8260
 QLast Update : Wed Aug 16 02:42:02 2023
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.230	168	291736	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	9.106	114	531571	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.871	117	516871	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.794	152	202281	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.583	65	205150	45.191	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	90.380%
35) Dibromofluoromethane	8.171	113	180779	50.522	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.040%
50) Toluene-d8	10.571	98	644085	48.477	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	96.960%
62) 4-Bromofluorobenzene	12.853	95	238373	54.889	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	109.780%
Target Compounds						
					Qvalue	
16) Acetone	4.436	43	29866	15.944	ug/l	90
20) Methylene Chloride	5.277	84	18385	4.606	ug/l #	77
25) 2-Butanone	7.500	43	28174	9.382	ug/l #	79
40) Benzene	8.612	78	414534	26.055	ug/l	98
43) Isopropyl Acetate	8.694	43	239556	22.505	ug/l #	79
51) 4-Methyl-2-Pentanone	10.447	43	87793	13.906	ug/l	89
52) Toluene	10.635	92	316580	32.410	ug/l	97
67) Ethyl Benzene	11.965	91	50241	2.511	ug/l	91
68) m/p-Xylenes	12.071	106	89576	11.974	ug/l	95
69) o-Xylene	12.400	106	57168	7.856	ug/l	93
84) 1,2,4-Trimethylbenzene	13.488	105	39783	3.006	ug/l	93
95) Naphthalene	15.647	128	334574	36.619	ug/l	100

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

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