

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN091522\
 Data File : VN074469.D
 Acq On : 15 Sep 2022 18:22
 Operator : JC\MD
 Sample : N4578-08
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 CHERRY-HILL-COMP-2

Quant Time: Sep 15 19:28:28 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N091522W.M
 Quant Title : SW846 8260
 QLast Update : Thu Sep 15 19:20:33 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	8.016	168	438540	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.904	114	868872	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.686	117	893140	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.615	152	400143	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.369	65	326242	43.845	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	87.680%
35) Dibromofluoromethane	7.951	113	269034	46.612	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.220%
50) Toluene-d8	10.380	98	1029948	46.141	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	92.280%
62) 4-Bromofluorobenzene	12.674	95	424088	55.472	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	110.940%
Target Compounds						
					Qvalue	
16) Acetone	4.222	43	81536	17.124	ug/l	95
20) Methylene Chloride	5.004	84	25225	3.499	ug/l #	91
37) Ethyl Acetate	7.339	43	114706	7.270	ug/l	96
43) Isopropyl Acetate	8.498	43	560783	22.691	ug/l #	83
68) m/p-Xylenes	11.904	106	4800	0.315	ug/l	98

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

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