

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX021425\
 Data File : VX044962.D
 Acq On : 14 Feb 2025 15:29
 Operator : JC/MD
 Sample : Q1356-11
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 CARBON-SOLID

Quant Time: Feb 14 22:15:29 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X021025W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 11 03:41:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	78767	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	161354	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	147582	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	65157	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	64334	55.797	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	111.600%
35) Dibromofluoromethane	5.379	113	52974	50.496	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	101.000%
50) Toluene-d8	8.647	98	199003	50.165	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	100.340%
62) 4-Bromofluorobenzene	11.079	95	69016	51.607	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	=	103.220%
Target Compounds						
						Qvalue
16) Acetone	2.380	43	14951	32.255	ug/l	96
18) Methyl Acetate	2.709	43	2102	1.437	ug/l	91
30) Chloroform	5.099	83	2631	1.397	ug/l	95
43) Isopropyl Acetate	6.342	43	22497	8.275	ug/l	98

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

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