

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042225\
 Data File : VX045892.D
 Acq On : 22 Apr 2025 18:00
 Operator : JC/MD
 Sample : Q1842-24
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PL-714-COMP-11

Quant Time: Apr 23 01:36:16 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Apr 02 03:11:43 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.550	168	65801	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	131202	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	125687	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	54408	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	64724	53.787	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 107.580%	
35) Dibromofluoromethane	5.379	113	46969	50.457	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 100.920%	
50) Toluene-d8	8.647	98	166402	51.214	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.420%	
62) 4-Bromofluorobenzene	11.079	95	63616	53.753	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 107.500%	
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	23810	48.187	ug/l	100
18) Methyl Acetate	2.709	43	2555	2.246	ug/l	92
20) Methylene Chloride	2.782	84	30975	33.485	ug/l	97
43) Isopropyl Acetate	6.342	43	10494	4.371	ug/l	98

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX042225\
Data File : VX045892.D
Acq On : 22 Apr 2025 18:00
Operator : JC/MD
Sample : Q1842-24
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 19 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
PL-714-COMP-11

Quant Time: Apr 23 01:36:16 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X040225W.M
Quant Title : SW846 8260
QLast Update : Wed Apr 02 03:11:43 2025
Response via : Initial Calibration

