

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020524\
 Data File : VX040115.D
 Acq On : 05 Feb 2024 14:54
 Operator : JC/MD
 Sample : P1141-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 EB-3-2-4

Quant Time: Feb 06 03:01:24 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X013024W.M
 Quant Title : SW846 8260
 QLast Update : Thu Feb 01 05:44:39 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	95695	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	173310	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	164983	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	56590	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	68618	44.555	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	89.100%
35) Dibromofluoromethane	5.385	113	55461	47.010	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.020%
50) Toluene-d8	8.647	98	212637	48.195	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	96.380%
62) 4-Bromofluorobenzene	11.079	95	62065	36.533	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	73.060%#
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	19828	27.478	ug/l #	89
20) Methylene Chloride	2.788	84	3943	3.075	ug/l	97
37) Ethyl Acetate	4.733	43	3317	1.780	ug/l #	92
43) Isopropyl Acetate	6.336	43	77152	26.255	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020524\
Data File : VX040115.D
Acq On : 05 Feb 2024 14:54
Operator : JC/MD
Sample : P1141-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 16 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
EB-3-2-4

Quant Time: Feb 06 03:01:24 2024
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X013024W.M
Quant Title : SW846 8260
QLast Update : Thu Feb 01 05:44:39 2024
Response via : Initial Calibration

