

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020524\
 Data File : VX040122.D
 Acq On : 05 Feb 2024 17:35
 Operator : JC/MD
 Sample : P1337-06
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 TP-2

Quant Time: Feb 06 03:04:07 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	91074	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	162850	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	160867	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	74475	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	65704	44.827	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	89.660%
35) Dibromofluoromethane	5.391	113	51631	46.574	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	93.140%
50) Toluene-d8	8.647	98	202807	48.919	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	97.840%
62) 4-Bromofluorobenzene	11.079	95	81230	50.885	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	101.780%
Target Compounds						
					Qvalue	
16) Acetone	2.380	43	32827	47.800	ug/l	95
20) Methylene Chloride	2.788	84	5763	4.722	ug/l	93
37) Ethyl Acetate	4.733	43	3133	1.789	ug/l #	90
43) Isopropyl Acetate	6.342	43	74296	26.907	ug/l	99

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX020524\
Data File : VX040122.D
Acq On : 05 Feb 2024 17:35
Operator : JC/MD
Sample : P1337-06
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 23 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
TP-2

Quant Time: Feb 06 03:04:07 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

