

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090722\
 Data File : VX031173.D
 Acq On : 07 Sep 2022 20:46
 Operator : JC/MD
 Sample : N4357-10DL 20X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 DUP07-20220823DL

Quant Time: Sep 07 23:27:13 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090722W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 07 23:22:41 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	78762	50.000	ug/l	# 0.00
34) 1,4-Difluorobenzene	6.763	114	141467	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	123901	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	56137	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	105516	56.585	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	=	113.180%
35) Dibromofluoromethane	5.391	113	80008	51.899	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	103.800%
50) Toluene-d8	8.653	98	277662	49.574	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	=	99.140%
62) 4-Bromofluorobenzene	11.079	95	107631	52.437	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	104.880%
Target Compounds						
						Qvalue
9) 1,1,2-Trichlorotrifluo...	2.331	101	785	0.555	ug/l	91
44) Trichloroethene	7.129	130	56663	32.142	ug/l	96

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090722\
Data File : VX031173.D
Acq On : 07 Sep 2022 20:46
Operator : JC/MD
Sample : N4357-10DL 20X
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 26 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
DUP07-20220823DL

Quant Time: Sep 07 23:27:13 2022
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090722W.M
Quant Title : SW846 8260
QLast Update : Wed Sep 07 23:22:41 2022
Response via : Initial Calibration

