

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112425\
 Data File : VX048664.D
 Acq On : 24 Nov 2025 15:27
 Operator : JC/MD
 Sample : Q3707-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 V933

Quant Time: Nov 25 17:46:28 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
 Quant Title : SW846 8260
 QLast Update : Sat Nov 08 05:08:11 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	177220	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.745	114	315823	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.037	117	342865	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.006	152	206089	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.946	65	130755	52.950	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	105.900%
35) Dibromofluoromethane	5.379	113	112434	47.037	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	94.080%
50) Toluene-d8	8.634	98	365310	49.001	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	98.000%
62) 4-Bromofluorobenzene	11.061	95	170710	61.445	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	122.880%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.940	59	6926	14.801	ug/l	96
16) Acetone	2.373	43	51967	44.458	ug/l	96
20) Methylene Chloride	2.794	84	6470	2.692	ug/l #	86

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

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX112425\
Data File : VX048664.D
Acq On : 24 Nov 2025 15:27
Operator : JC/MD
Sample : Q3707-03
Misc : 5.0mL/MSVOA_X/WATER
ALS Vial : 8 Sample Multiplier: 1

Instrument :
MSVOA_X
ClientSampleId :
V933

Quant Time: Nov 25 17:46:28 2025
Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X110725W.M
Quant Title : SW846 8260
QLast Update : Sat Nov 08 05:08:11 2025
Response via : Initial Calibration

