

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX031125\
 Data File : VX045230.D
 Acq On : 11 Mar 2025 17:29
 Operator : JC/MD
 Sample : Q1531-19
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 21 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 RE104D2-20250307

Quant Time: Mar 12 01:52:39 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022825W.M
 Quant Title : SW846 8260
 QLast Update : Fri Feb 28 06:45:16 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.544	168	68480	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.757	114	133865	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.049	117	123884	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.018	152	53817	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.952	65	58163	53.396	ug/l	0.00
Spiked Amount	50.000	Range	74 - 125	Recovery	= 106.800%	
35) Dibromofluoromethane	5.379	113	46770	52.250	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 104.500%	
50) Toluene-d8	8.647	98	166263	51.235	ug/l	0.00
Spiked Amount	50.000	Range	86 - 113	Recovery	= 102.460%	
62) 4-Bromofluorobenzene	11.079	95	59299	55.145	ug/l	0.00
Spiked Amount	50.000	Range	77 - 121	Recovery	= 110.280%	
Target Compounds						
					Qvalue	
12) 1,1-Dichloroethene	2.307	96	1494	1.775	ug/l	87
16) Acetone	2.386	43	832	1.646	ug/l #	85
27) cis-1,2-Dichloroethene	4.483	96	49010	47.757	ug/l	95
30) Chloroform	5.099	83	3108	1.832	ug/l	95
42) 1,2-Dichloroethane	6.086	62	2436	1.745	ug/l	85
44) Trichloroethene	7.123	130	396610	435.736	ug/l	99
64) Tetrachloroethene	9.275	164	471	0.594	ug/l #	85

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

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