

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092525\
 Data File : VV039177.D
 Acq On : 25 Sep 2025 13:40
 Operator : SY/MD
 Sample : Q3168-05
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 TOTES COMP#1

Quant Time: Sep 26 04:36:29 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\82V092225W.M
 Quant Title : SW846 8260
 QLast Update : Wed Sep 24 08:08:08 2025
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	4.603	168	424278	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.535	114	802353	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	790092	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.175	152	406442	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.944	65	405843	66.092	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	132.180%#
35) Dibromofluoromethane	4.503	113	363530	56.359	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	112.720%
50) Toluene-d8	7.239	98	866364	45.732	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	91.460%
62) 4-Bromofluorobenzene	10.001	95	374455	48.014	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	96.020%
Target Compounds						
						Qvalue
11) Tert butyl alcohol	2.574	59	28626	22.057	ug/l #	88
16) Acetone	2.121	43	439192	109.225	ug/l	99
17) Carbon Disulfide	2.256	76	21890	1.119	ug/l	99
25) 2-Butanone	3.883	43	165528	33.663	ug/l	100
51) 4-Methyl-2-Pentanone	7.165	43	28490	2.684	ug/l	92
52) Toluene	7.326	92	5165	0.281	ug/l	90
59) 2-Hexanone	8.079	43	97296	13.576	ug/l	95
68) m/p-Xylenes	9.075	106	7809	0.607	ug/l	83
69) o-Xylene	9.474	106	3531	0.310	ug/l	92

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

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