

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\DATA\VV092525\
 Data File : VV039183.D
 Acq On : 25 Sep 2025 15:55
 Operator : SY/MD
 Sample : Q3168-05RE
 Misc : 5.0mL/MSVOA_V/WATER
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 TOTES COMP#1RE

Quant Time: Sep 26 01:31:22 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	4.600	168	518499	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	5.535	114	966445	50.000	ug/l	0.00
63) Chlorobenzene-d5	8.776	117	862048	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	11.172	152	399824	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	4.944	65	407460	54.297	ug/l	0.00
Spiked Amount	50.000	Range	81 - 118	Recovery	=	108.600%
35) Dibromofluoromethane	4.503	113	364495	46.914	ug/l	0.00
Spiked Amount	50.000	Range	80 - 119	Recovery	=	93.820%
50) Toluene-d8	7.239	98	887221	38.881	ug/l	0.00
Spiked Amount	50.000	Range	89 - 112	Recovery	=	77.760%#
62) 4-Bromofluorobenzene	10.001	95	366650	39.031	ug/l	0.00
Spiked Amount	50.000	Range	85 - 114	Recovery	=	78.060%#
Target Compounds						
					Qvalue	
11) Tert butyl alcohol	2.577	59	31850	20.081	ug/l	99
16) Acetone	2.121	43	453120	91.092	ug/l	100
17) Carbon Disulfide	2.256	76	20156	0.843	ug/l #	93
25) 2-Butanone	3.886	43	177245	29.496	ug/l	95
51) 4-Methyl-2-Pentanone	7.162	43	36643	2.866	ug/l	98
52) Toluene	7.333	92	6169	0.278	ug/l	100
59) 2-Hexanone	8.075	43	108220	12.714	ug/l	98
68) m/p-Xylenes	9.072	106	7137	0.509	ug/l	98
69) o-Xylene	9.480	106	3828	0.308	ug/l	95

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

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