

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX022924\
 Data File : VX040538.D
 Acq On : 29 Feb 2024 15:36
 Operator : JC/MD
 Sample : P1641-05
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 OF-3

Quant Time: Mar 01 03:48:54 2024
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X022624W.M
 Quant Title : SW846 8260
 QLast Update : Tue Feb 27 05:01:33 2024
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	70286	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	124256	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	108839	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	62135	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	58929	50.128	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	= 100.260%	
35) Dibromofluoromethane	5.385	113	41526	48.059	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	= 96.120%	
50) Toluene-d8	8.647	98	153476	48.933	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	= 97.860%	
62) 4-Bromofluorobenzene	11.079	95	62294	52.103	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	= 104.200%	
Target Compounds						
					Qvalue	
13) Acrolein	2.233	56	564	4.451	ug/l	91
16) Acetone	2.380	43	9478	24.723	ug/l	99
25) 2-Butanone	4.580	43	2095	3.136	ug/l	92
40) Benzene	6.037	78	1057	0.293	ug/l	97
51) 4-Methyl-2-Pentanone	8.580	43	6618	4.378	ug/l	99

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

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