

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090622\
 Data File : VX031138.D
 Acq On : 06 Sep 2022 19:10
 Operator : JC/MD
 Sample : N4489-03DL 10X
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ORA-0044DL

Quant Time: Sep 07 01:14:11 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X090122W.M
 Quant Title : SW846 8260
 QLast Update : Fri Sep 02 02:49:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.556	168	213678	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	335562	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	304447	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	150510	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	130547	47.482	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	94.960%		
35) Dibromofluoromethane	5.391	113	124397	50.799	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	101.600%		
50) Toluene-d8	8.653	98	435177	48.286	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	96.580%		
62) 4-Bromofluorobenzene	11.079	95	162325	50.994	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	101.980%		
Target Compounds						Qvalue
11) Tert butyl alcohol	2.977	59	20842	17.521	ug/l	98
16) Acetone	2.386	43	213140	200.811	ug/l	100
18) Methyl Acetate	2.709	43	1824	0.598	ug/l #	67
25) 2-Butanone	4.568	43	87636	53.482	ug/l #	84
37) Ethyl Acetate	4.733	43	10681	3.584	ug/l #	92
51) 4-Methyl-2-Pentanone	8.580	43	5005	1.672	ug/l	89
59) 2-Hexanone	9.433	43	10767	4.755	ug/l	89

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

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