

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX090622\
 Data File : VX031130.D
 Acq On : 06 Sep 2022 16:02
 Operator : JC/MD
 Sample : N4489-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 8 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 ORA-0044

Quant Time: Sep 06 17:30:02 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	227806	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.763	114	369249	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	312998	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	155340	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	138282	47.176	ug/l	0.00
Spiked Amount 50.000	Range 74 - 125		Recovery =	94.360%		
35) Dibromofluoromethane	5.391	113	130248	48.336	ug/l	0.00
Spiked Amount 50.000	Range 75 - 124		Recovery =	96.680%		
50) Toluene-d8	8.653	98	447146	45.087	ug/l	0.00
Spiked Amount 50.000	Range 86 - 113		Recovery =	90.180%		
62) 4-Bromofluorobenzene	11.085	95	159668	45.583	ug/l	0.00
Spiked Amount 50.000	Range 83 - 123		Recovery =	91.160%		
Target Compounds						Qvalue
11) Tert butyl alcohol	2.977	59	133238	222.609	ug/l	97
16) Acetone	2.386	43	1961572	1733.491	ug/l	99
18) Methyl Acetate	2.709	43	18557	5.707	ug/l #	84
25) 2-Butanone	4.562	43	830837	475.595	ug/l #	85
37) Ethyl Acetate	4.727	43	116221	35.439	ug/l #	93
51) 4-Methyl-2-Pentanone	8.574	43	43897	13.323	ug/l	88
52) Toluene	8.720	92	1867	0.266	ug/l	93
59) 2-Hexanone	9.433	43	106822	42.872	ug/l	91
68) m/p-Xylenes	10.305	106	2545	0.507	ug/l	93
69) o-Xylene	10.640	106	1432	0.291	ug/l	96

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

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