

Data Path : Z:\voasrv\HPCHEM1\MSVOA_X\Data\VX010522\
 Data File : VX026347.D
 Acq On : 06 Jan 2022 01:39
 Operator : JC/MD
 Sample : N1019-03
 Misc : 5.0mL/MSVOA_X/WATER
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 PZ-9

Quant Time: Jan 06 05:04:46 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_X\Method\82X123021W.M
 Quant Title : SW846 8260
 QLast Update : Thu Dec 30 08:02:21 2021
 Response via : Initial Calibration

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

Internal Standards						
1) Pentafluorobenzene	5.562	168	153805	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	6.769	114	259221	50.000	ug/l	0.00
63) Chlorobenzene-d5	10.055	117	225369	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.024	152	100380	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	5.958	65	110936	49.857	ug/l	0.00
Spiked Amount	50.000	Range	78 - 117	Recovery	=	99.720%
35) Dibromofluoromethane	5.391	113	82761	49.108	ug/l	0.00
Spiked Amount	50.000	Range	75 - 124	Recovery	=	98.220%
50) Toluene-d8	8.653	98	300652	47.962	ug/l	0.00
Spiked Amount	50.000	Range	92 - 112	Recovery	=	95.920%
62) 4-Bromofluorobenzene	11.079	95	113914	51.047	ug/l	0.00
Spiked Amount	50.000	Range	83 - 123	Recovery	=	102.100%
Target Compounds						
					Qvalue	
16) Acetone	2.386	43	7923	6.572	ug/l #	89
31) Cyclohexane	5.477	56	7341	2.317	ug/l	93
37) Ethyl Acetate	4.721	43	16065	5.097	ug/l	99
40) Benzene	6.044	78	497996	64.194	ug/l	99
52) Toluene	8.720	92	112895	23.639	ug/l	99
67) Ethyl Benzene	10.195	91	347438	39.248	ug/l	97
68) m/p-Xylenes	10.299	106	207801	62.869	ug/l	92
69) o-Xylene	10.640	106	20589	6.366	ug/l	92
73) Isopropylbenzene	10.964	105	14216	1.822	ug/l	95
78) n-propylbenzene	11.305	91	10071	1.123	ug/l	92
84) 1,2,4-Trimethylbenzene	11.756	105	21335	3.298	ug/l	97

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

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