

Data Path : Z:\VOASRV\HPCHEM1\MSVOA X\DATA\VX061120\
 Data File : VX016715.D
 Acq On : 11 Jun 2020 19:21
 Operator : JC/SP
 Sample : L2977-06
 Misc : 5.0mL/MSVOA X/WATER
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 MSVOA_X
 ClientSampleId :
 I-LDC-2

Quant Time: Jun 12 15:55:11 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_X\METHOD\82X052720W.M
 Quant Title : SW846 8260
 QLast Update : Wed May 27 16:31:05 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	5.64	168	209354	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	6.83	114	338826	50.00	ug/l	0.00
63) Chlorobenzene-d5	10.10	117	334012	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	12.06	152	177894	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	6.03	65	123091	47.34	ug/l	0.00
Spiked Amount			Recovery	=	94.68%	
35) Dibromofluoromethane	5.47	113	102943	49.34	ug/l	0.00
Spiked Amount			Recovery	=	98.68%	
50) Toluene-d8	8.70	98	407828	50.22	ug/l	0.00
Spiked Amount			Recovery	=	100.44%	
62) 4-Bromofluorobenzene	11.12	95	167509	53.39	ug/l	0.00
Spiked Amount			Recovery	=	106.78%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	2.42	43	477601	419.287	ug/l	100
17) Carbon Disulfide	2.55	76	2074292	346.748	ug/l	99
25) 2-Butanone	4.64	43	91074	47.186	ug/l	93
30) Chloroform	5.18	83	141848	35.047	ug/l	98
40) Benzene	6.12	78	174401	18.255	ug/l	98
47) Bromodichloromethane	7.88	83	8337	2.465	ug/l	94
52) Toluene	8.77	92	50470	8.378	ug/l	97
67) Ethyl Benzene	10.24	91	60046	4.953	ug/l	100
68) m/p-Xylenes	10.34	106	87216	19.402	ug/l	98
69) o-Xylene	10.68	106	39483	9.069	ug/l	96
78) n-propylbenzene	11.34	91	3723	0.270	ug/l	92
79) 2-Chlorotoluene	11.41	91	3541	0.421	ug/l	76
80) 1,3,5-Trimethylbenzene	11.49	105	16151	1.569	ug/l	97
84) 1,2,4-Trimethylbenzene	11.79	105	61758	5.933	ug/l	99
95) Naphthalene	13.82	128	6516323	527.446	ug/l	99

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

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