

Data Path : Z:\VOASRV\HPCHEM1\MSVOA N\DATA\VN032819\
 Data File : VN054747.D
 Acq On : 28 Mar 2019 16:25
 Operator : JC/SP
 Sample : K2112-01
 Misc : 5.00mL/MSVOA N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 EPH-4-8-WC-001

Quant Time: Mar 29 07:26:07 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N032019W.M
 Quant Title : SW846 8260
 QLast Update : Fri Mar 22 02:09:07 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	546719	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	879161	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	742035	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.34	152	279242	50.00	ug/l	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
33) 1,2-Dichloroethane-d4	8.03	65	350526	51.55	ug/l	0.00
Spiked Amount			Recovery	=	103.10%	
35) Dibromofluoromethane	7.59	113	275584	48.66	ug/l	0.00
Spiked Amount			Recovery	=	97.32%	
50) Toluene-d8	10.09	98	1066452	50.08	ug/l	0.00
Spiked Amount			Recovery	=	100.16%	
62) 4-Bromofluorobenzene	12.40	95	331130	45.70	ug/l	0.00
Spiked Amount			Recovery	=	91.40%	

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.82	43	14568	7.125	ug/l #	78
20) Methylene Chloride	4.55	84	2805534	465.508	ug/l #	83
24) 1,1-Dichloroethane	5.85	63	36246	3.437	ug/l	94
27) cis-1,2-Dichloroethene	6.83	96	34938	5.440	ug/l	100
30) Chloroform	7.37	83	45667	4.437	ug/l	97
40) Benzene	8.04	78	195774	8.329	ug/l	100
42) 1,2-Dichloroethane	8.12	62	16490	1.926	ug/l	99
43) Isopropyl Acetate	8.17	43	20384	2.064	ug/l #	78
44) Trichloroethene	8.83	130	120145	18.384	ug/l	87
52) Toluene	10.16	92	21914	1.516	ug/l	91
64) Tetrachloroethene	10.63	164	49951	8.116	ug/l	96
68) m/p-Xylenes	11.62	106	10472	1.037	ug/l	96
73) Isopropylbenzene	12.25	105	63924	2.841	ug/l	98
78) n-propylbenzene	12.59	91	34748	1.424	ug/l	96
85) sec-Butylbenzene	13.17	105	48120	2.405	ug/l	85

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

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