

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031419\
 Data File : VN054264.D
 Acq On : 14 Mar 2019 16:14
 Operator : JC/SP
 Sample : K1899-01
 Misc : 5.00mL/MSVOA_N/WATER
 ALS Vial : 13 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 MW-1A

Manual Integrations
 APPROVED

MMDadoda

Quant Time: Mar 16 04:20:13 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
 Quant Title : SW846 8260
 QLast Update : Wed Mar 13 03:33:31 2019
 Response via : Initial Calibration

Internal Standards R.T. QIon Response Conc Units Dev(Min) 3/18/2019 10:51:24 AM

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) Pentafluorobenzene	7.67	168	1379221	50.00	ug/l	0.00
34) 1,4-Difluorobenzene	8.59	114	2215978	50.00	ug/l	0.00
63) Chlorobenzene-d5	11.41	117	2441287	50.00	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.35	152	1349842	50.00	ug/l	0.00

System Monitoring Compounds

33) 1,2-Dichloroethane-d4	8.03	65	705231	48.15	ug/l	0.00
Spiked Amount	50.000		Recovery	=	96.30%	
35) Dibromofluoromethane	7.59	113	674080	47.16	ug/l	0.00
Spiked Amount	50.000		Recovery	=	94.32%	
50) Toluene-d8	10.09	98	2833114	53.67	ug/l	0.00
Spiked Amount	50.000		Recovery	=	107.34%	
62) 4-Bromofluorobenzene	12.40	95	1204636	65.23	ug/l	0.00
Spiked Amount	50.000		Recovery	=	130.46%	

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
16) Acetone	3.82	43	418903	99.884	ug/l	99
25) 2-Butanone	6.85	43	111752	22.703	ug/l	96
29) Tetrahydrofuran	7.23	42	6783	2.143	ug/l #	81
30) Chloroform	7.37	83	93887	3.819	ug/l	97
39) Methylcyclohexane	9.08	83	10961	0.437	ug/l	87
40) Benzene	8.04	78	19458501	333.208	ug/l	100
42) 1,2-Dichloroethane	8.13	62	32155	1.631	ug/l	97
51) 4-Methyl-2-Pentanone	9.99	43	59919	5.501	ug/l	97
52) Toluene	10.15	92	17058884m	460.704	ug/l	
67) Ethyl Benzene	11.50	91	17034781m	191.708	ug/l	
68) m/p-Xylenes	11.62	106	14792364	425.928	ug/l #	1
69) o-Xylene	11.95	106	8309877	245.450	ug/l	89
73) Isopropylbenzene	12.25	105	1941785	17.980	ug/l	100
78) n-propylbenzene	12.59	91	1035108	8.833	ug/l	99
80) 1,3,5-Trimethylbenzene	12.73	105	4486124	49.940	ug/l	99
84) 1,2,4-Trimethylbenzene	13.04	105	15523713	176.151	ug/l	95
85) sec-Butylbenzene	13.17	105	118627m	1.106	ug/l	
86) p-Isopropyltoluene	13.29	119	4213013	44.785	ug/l	98
87) 1,3-Dichlorobenzene	13.28	146	12807m	0.266	ug/l	
88) 1,4-Dichlorobenzene	13.36	146	270317	5.714	ug/l	97
89) n-Butylbenzene	13.62	91	104631m	1.464	ug/l	
91) 1,2-Dichlorobenzene	13.65	146	139081	2.954	ug/l #	81
95) Naphthalene	15.11	128	37259305m	589.132	ug/l	

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA_N\DATA\VN031419\
Data File : VN054264.D
Acq On : 14 Mar 2019 16:14
Operator : JC/SP
Sample : K1899-01
Misc : 5.00mL/MSVOA_N/WATER
ALS Vial : 13 Sample Multiplier: 1

Instrument :
MSVOA_N
ClientSampleId :
MW-1A

Manual Integrations
APPROVED

MMDadoda

Quant Time: Mar 16 04:20:13 2019
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_N\METHODS\82N031219W.M
Quant Title : SW846 8260
QLast Update : Wed Mar 13 03:33:31 2019
Response via : Initial Calibration

