

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\U110118\
 Data File : VU027900.D
 Acq On : 01 Nov 2018 13:35
 Operator : MD/SY
 Sample : MDL07
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL07

Manual Integrations
 APPROVED

MMDadoda
 11/5/2018 4:10:26 PM

Quant Time: Nov 02 07:46:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR103018WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Fri Nov 02 08:02:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.89	114	78795	5.00	ug/L	0.00
28) Chlorobenzene-d5	9.09	117	77130	5.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	11.48	152	37938	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.40	65	23663	3.84	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	76.80%
7) Chloroethane-d5	1.68	69	19479	4.10	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	82.00%
11) 1,1-Dichloroethene-d2	2.28	63	38203	3.27	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	65.40%
20) 2-Butanone-d5	4.20	46	86978	47.36	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	94.72%
24) Chloroform-d	4.65	84	45333	4.09	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	81.80%
26) 1,2-Dichloroethane-d4	5.31	65	30684	4.82	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	96.40%
32) Benzene-d6	5.34	84	88931	4.21	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	84.20%
36) 1,2-Dichloropropane-d6	6.33	67	31914	4.27	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	85.40%
41) Toluene-d8	7.57	98	81840	4.07	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	81.40%
43) trans-1,3-Dichloropropene-	7.85	79	11744	4.06	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	81.20%
46) 2-Hexanone-d5	8.31	63	68796	43.46	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	86.92%
57) 1,1,2,2-Tetrachloroethane-	10.43	84	23194	4.38	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	87.60%
64) 1,2-Dichlorobenzene-d4	11.86	152	31479	4.34	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	86.80%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.21	85	2463	0.318 ug/L	95
3) Chloromethane	1.33	50	2261	0.294 ug/L	96
5) Vinyl chloride	1.40	62	2511	0.328 ug/L	87
6) Bromomethane	1.63	94	821	0.326 ug/L	89
8) Chloroethane	1.70	64	1719	0.383 ug/L #	82
9) Trichlorofluoromethane	1.89	101	3259	0.320 ug/L	89
10) 1,1,2-Trichloro-1,2,2-trif	2.29	101	1714	0.330 ug/L #	85
12) 1,1-Dichloroethene	2.29	96	1520	0.311 ug/L #	1
13) Acetone	2.32	43	4245	3.667 ug/L	88
14) Carbon disulfide	2.48	76	4311	0.247 ug/L	98
15) Methyl Acetate	2.62	43	1234m	0.398 ug/L	
16) Methylene chloride	2.70	84	1919	0.349 ug/L	87
17) Methyl tert-butyl Ether	3.00	73	3817	0.262 ug/L #	77
18) trans-1,2-Dichloroethene	2.99	96	1547	0.297 ug/L	88
19) 1,1-Dichloroethane	3.45	63	3168	0.288 ug/L	96
21) 2-Butanone	4.29	43	6125	3.043 ug/L	89
22) cis-1,2-Dichloroethene	4.22	96	1625	0.270 ug/L	73

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110118\
 Data File : VU027900.D
 Acq On : 01 Nov 2018 13:35
 Operator : MD/SY
 Sample : MDL07
 Misc : 25.0mL/MSVOA U/WATER
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 MDL07

Manual Integrations
 APPROVED

MMDadoda
 11/5/2018 4:10:26 PM

Quant Time: Nov 02 07:46:56 2018
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR103018WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Fri Nov 02 08:02:52 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)

23) Bromochloromethane	4.55	128	671	0.269	ug/L #	74
25) Chloroform	4.68	83	7525	0.676	ug/L	98
27) 1,2-Dichloroethane	5.41	62	2257	0.290	ug/L #	78
29) 1,1,1-Trichloroethane	4.92	97	2768	0.278	ug/L	95
30) Cyclohexane	4.99	56	3172	0.271	ug/L	96
31) Carbon tetrachloride	5.13	117	2439	0.286	ug/L	91
33) Benzene	5.40	78	6837	0.280	ug/L	100
34) Trichloroethene	6.19	95	1785	0.281	ug/L	93
35) Methylcyclohexane	6.41	83	3054	0.282	ug/L	92
37) 1,2-Dichloropropane	6.43	63	1893	0.272	ug/L #	90
38) Bromodichloromethane	6.77	83	2164	0.266	ug/L #	92
39) cis-1,3-Dichloropropene	7.27	75	2698	0.271	ug/L	99
40) 4-Methyl-2-pentanone	7.47	43	13875	2.822	ug/L #	89
42) Toluene	7.64	91	7435	0.291	ug/L	97
44) trans-1,3-Dichloropropene	7.88	75	2320	0.271	ug/L	99
45) 1,1,2-Trichloroethane	8.08	97	1183	0.276	ug/L	88
47) Tetrachloroethene	8.23	164	1457	0.298	ug/L	83
48) 2-Hexanone	8.37	43	11102	3.133	ug/L	96
49) Dibromochloromethane	8.48	129	1276	0.247	ug/L	88
50) 1,2-Dibromoethane	8.59	107	1044	0.255	ug/L #	74
51) Chlorobenzene	9.12	112	4986	0.315	ug/L	86
52) Ethylbenzene	9.25	91	7982	0.274	ug/L	88
53) m,p-xylene	9.38	106	2880	0.273	ug/L	84
54) o-xylene	9.78	106	2675	0.258	ug/L	74
55) Styrene	9.79	104	4653	0.267	ug/L	95
56) Isopropylbenzene	10.17	105	7803	0.276	ug/L	92
58) 1,1,2,2-Tetrachloroethane	10.46	83	1668	0.299	ug/L #	94
59) 1,2,3-Trichloropropane	10.49	75	1027m	0.244	ug/L	
61) Bromoform	9.96	173	639	0.229	ug/L #	74
62) 1,3-Dichlorobenzene	11.42	146	3482	0.285	ug/L	95
63) 1,4-Dichlorobenzene	11.51	146	3573	0.293	ug/L	95
65) 1,2-Dichlorobenzene	11.88	146	3232	0.281	ug/L	94
66) 1,2-Dibromo-3-chloropropan	12.65	75	182	0.202	ug/L #	42
67) 1,3,5-Trichlorobenzene	12.88	180	2701	0.275	ug/L	92
68) 1,2,4-trichlorobenzene	13.51	180	2234	0.268	ug/L #	91
69) Naphthalene	13.75	128	2451	0.189	ug/L #	89
70) 1,2,3-Trichlorobenzene	13.99	180	1949	0.256	ug/L	96

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

Data Path : Z:\VOASRV\HPCHEM1\MSVOA U\DATA\VU110118\
Data File : VU027900.D
Acq On : 01 Nov 2018 13:35
Operator : MD/SY
Sample : MDL07
Misc : 25.0mL/MSVOA U/WATER
ALS Vial : 11 Sample Multiplier: 1

Instrument :
MSVOA_U
ClientSampleId :
MDL07

Manual Integrations
APPROVED

MMDadoda
11/5/2018 4:10:26 PM

Quant Time: Nov 02 07:46:56 2018
Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_U\METHOD\SOMUTR103018WMA.M
Quant Title : TRACE VOA SOM01.0
QLast Update : Fri Nov 02 08:02:52 2018
Response via : Initial Calibration

