

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY111319\
 Data File : VY000670.D
 Acq On : 13 Nov 2019 14:04
 Operator : SY/MD
 Sample : MDL07
 Misc : 5.00G/10ML/MSVOA Y/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 MDL07

Manual Integrations
 APPROVED

MMDADODA
 11/19/2019 11:59:50 AM

Quant Time: Nov 14 08:44:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\SOM2YLM111219S.M
 Quant Title : VOC Analysis
 QLast Update : Thu Nov 14 08:38:56 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	8.70	114	338889	25.00	ug/L	0.00
28) Chlorobenzene-d5	11.49	117	306523	25.00	ug/L	0.00
60) 1,4-Dichlorobenzene-d4	13.42	152	140911	25.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	2.24	65	104962	21.93	ug/L	0.00
Spiked Amount	25.000	Range	30 - 150	Recovery	=	87.72%
7) Chloroethane-d5	2.76	69	98414	23.14	ug/L	0.00
Spiked Amount	25.000	Range	30 - 150	Recovery	=	92.56%
10) 1,1-Dichloroethene-d2	3.85	63	145265	16.43	ug/L	0.00
Spiked Amount	25.000	Range	45 - 110	Recovery	=	65.72%
20) 2-Butanone-d5	6.91	46	78410	39.71	ug/L	0.00
Spiked Amount	50.000	Range	20 - 135	Recovery	=	79.42%
24) Chloroform-d	7.49	84	193920	21.81	ug/L	0.00
Spiked Amount	25.000	Range	40 - 150	Recovery	=	87.24%
26) 1,2-Dichloroethane-d4	8.15	65	113311	21.98	ug/L	0.00
Spiked Amount	25.000	Range	70 - 130	Recovery	=	87.92%
29) Benzene-d6	8.12	84	409804	23.02	ug/L	0.00
Spiked Amount	25.000	Range	20 - 135	Recovery	=	92.08%
33) 1,2-Dichloropropane-d6	9.13	67	126073	22.60	ug/L	0.00
Spiked Amount	25.000	Range	70 - 120	Recovery	=	90.40%
37) Toluene-d8	10.18	98	370896	22.30	ug/L	0.00
Spiked Amount	25.000	Range	30 - 130	Recovery	=	89.20%
38) trans-1,3-Dichloropropene-	10.44	79	55750	21.42	ug/L	0.00
Spiked Amount	25.000	Range	30 - 135	Recovery	=	85.68%
39) 2-Hexanone-d5	10.79	63	60403	40.06	ug/L	0.00
Spiked Amount	50.000	Range	20 - 135	Recovery	=	80.12%
48) 1,1,2,2-Tetrachloroethane-	12.56	84	112612	20.75	ug/L	0.00
Spiked Amount	25.000	Range	45 - 120	Recovery	=	83.00%
61) 1,2-Dichlorobenzene-d4	13.72	152	131091	23.26	ug/L	0.00
Spiked Amount	25.000	Range	75 - 120	Recovery	=	93.04%

Target Compounds

					Ovalue
2) Dichlorodifluoromethane	1.90	85	2605	0.582 ug/L	88
3) Chloromethane	2.12	50	3022	0.538 ug/L	93
5) Vinyl chloride	2.25	62	3023	0.540 ug/L #	34
6) Bromomethane	2.66	94	1876m	0.613 ug/L	
8) Chloroethane	2.79	64	1877m	0.534 ug/L	
9) Trichlorofluoromethane	3.13	101	5040	0.635 ug/L #	63
11) 1,1,2-Trichloro-1,2,2-trif	3.91	101	3021m	0.671 ug/L	
12) 1,1-Dichloroethene	3.86	96	3710m	0.844 ug/L	
13) Acetone	3.98	43	3167m	2.192 ug/L	
14) Carbon disulfide	4.20	76	5908m	0.436 ug/L	
15) Methyl Acetate	4.49	43	1255m	0.379 ug/L	
16) Methylene chloride	4.72	84	13618	2.434 ug/L	83
17) Methyl tert-butyl Ether	5.24	73	6459m	0.483 ug/L	
18) trans-1,2-Dichloroethene	5.23	96	2279	0.464 ug/L	85
19) 1,1-Dichloroethane	6.02	63	4215	0.488 ug/L	93
22) cis-1,2-Dichloroethene	6.99	96	2794m	0.515 ug/L	
23) Bromochloromethane	7.34	128	1284m	0.542 ug/L	

Data Path : Z:\VOASRV\HPCHEM1\MSVOA Y\DATA\VY111319\
 Data File : VY000670.D
 Acq On : 13 Nov 2019 14:04
 Operator : SY/MD
 Sample : MDL07
 Misc : 5.00G/10ML/MSVOA Y/SOIL
 ALS Vial : 1 Sample Multiplier: 1

Instrument :
 MSVOA_Y
 ClientSampleId :
 MDL07

Manual Integrations
 APPROVED

MMDADODA
 11/19/2019 11:59:50 AM

Quant Time: Nov 14 08:44:30 2019
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_Y\METHODS\SOM2YLM111219S.M
 Quant Title : VOC Analysis
 QLast Update : Thu Nov 14 08:38:56 2019
 Response via : Initial Calibration

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

25) Chloroform	7.52	83	10970	1.179	ug/L #	70
27) 1,2-Dichloroethane	8.25	62	3429	0.551	ug/L	96
31) 1,1,1-Trichloroethane	7.71	97	3869	0.512	ug/L #	77
32) Carbon tetrachloride	7.91	117	3060m	0.465	ug/L	
34) Benzene	8.18	78	10578	0.528	ug/L	100
35) Trichloroethene	8.94	95	2919	0.556	ug/L	97
36) Methylcyclohexane	9.18	83	3733	0.415	ug/L #	83
40) 1,2-Dichloropropane	9.23	63	2268	0.447	ug/L	100
41) Bromodichloromethane	9.50	83	3673	0.562	ug/L	96
42) cis-1,3-Dichloropropene	9.94	75	3830	0.468	ug/L	97
43) 4-Methyl-2-pentanone	10.07	43	4594	0.946	ug/L #	53
44) Toluene	10.25	91	11165	0.515	ug/L	91
45) trans-1,3-Dichloropropene	10.47	75	2377	0.338	ug/L #	80
46) 1,1,2-Trichloroethane	10.64	97	2255	0.541	ug/L	88
47) Tetrachloroethene	10.72	164	2299	0.586	ug/L	80
49) 2-Hexanone	10.83	43	4506	1.191	ug/L #	94
50) Dibromochloromethane	10.99	129	2126	0.452	ug/L	96
51) 1,2-Dibromoethane	11.08	107	2061	0.508	ug/L #	78
52) Chlorobenzene	11.51	112	6201	0.451	ug/L	93
53) Ethylbenzene	11.59	91	11686	0.477	ug/L	95
54) m,p-Xylene	11.69	106	4857	0.524	ug/L	80
55) o-xylene	12.03	106	4700	0.516	ug/L	82
56) Styrene	12.04	104	6831	0.452	ug/L	90
57) Isopropylbenzene	12.32	105	10629	0.449	ug/L	91
58) 1,1,2,2-Tetrachloroethane	12.58	83	3115	0.599	ug/L #	81
59) 1,2,3-Trichloropropane	12.63	75	1868	0.442	ug/L	100
62) Bromoform	12.20	173	1128	0.421	ug/L #	90
63) 1,3-Dichlorobenzene	13.36	146	5147	0.512	ug/L #	81
64) 1,4-Dichlorobenzene	13.44	146	5575	0.555	ug/L #	87
65) 1,2-Dichlorobenzene	13.74	146	5289	0.568	ug/L	94
66) 1,2-Dibromo-3-chloropropan	14.36	75	488m	0.552	ug/L	
67) 1,3,5-Trichlorobenzene	14.50	180	3475m	0.496	ug/L	
68) 1,2,4-trichlorobenzene	15.00	180	3124	0.499	ug/L	93
69) Naphthalene	15.23	128	7178	0.460	ug/L #	93
70) 1,2,3-Trichlorobenzene	15.42	180	2804	0.482	ug/L	96

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

