

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082721\
 Data File : VU044532.D
 Acq On : 27 Aug 2021 15:43
 Operator : SY/MD
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD005097

Manual Integrations
 APPROVED

MMDadoda
 8/30/2021 10:54:22 AM

Quant Time: Aug 28 01:21:21 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR081821WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Sat Aug 28 01:19:29 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Difluorobenzene	6.256	114	124449	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.423	117	126482	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.819	152	64308	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.601	65	33703	3.349	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	67.000%
7) Chloroethane-d5	1.919	69	33754	3.859	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	77.200%
11) 1,1-Dichloroethene-d2	2.572	65	16357	3.820	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	76.400%
20) 2-Butanone-d5	4.639	46	139570m	50.140	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	100.280%
24) Chloroform-d	5.073	84	81773	4.595	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	91.800%
26) 1,2-Dichloroethane-d4	5.710	65	43442	4.476	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	89.600%
32) Benzene-d6	5.735	84	155864	4.376	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	87.600%
36) 1,2-Dichloropropane-d6	6.700	67	50703	4.450	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	89.000%
41) Toluene-d8	7.906	98	139465	4.416	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	88.400%
43) trans-1,3-Dichloroprop...	8.185	79	18828	4.601	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	92.000%
46) 2-Hexanone-d5	8.645	63	113086	51.105	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	102.200%
56) 1,1,2,2-Tetrachloroeth...	10.764	84	45814	4.774	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	95.400%
66) 1,2-Dichlorobenzene-d4	12.201	152	49540	4.560	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	91.200%
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.388	85	63994	5.000	ug/L	100
3) Chloromethane	1.520	50	80297	4.789	ug/L	100
5) Vinyl chloride	1.607	62	81262	5.243	ug/L	100
6) Bromomethane	1.861	94	44655	5.300	ug/L	100
8) Chloroethane	1.938	64	49188	5.191	ug/L	96
9) Trichlorofluoromethane	2.144	101	94533	5.334	ug/L	99
10) 1,1,2-Trichloro-1,2,2-...	2.588	101	51786	5.311	ug/L	99
12) 1,1-Dichloroethene	2.584	96	51362	5.370	ug/L	97
13) Acetone	2.646	43	106283m	45.981	ug/L	
14) Carbon disulfide	2.800	76	161911	5.219	ug/L	99
15) Methyl Acetate	2.961	43	17677	4.682	ug/L #	89
16) Methylene chloride	3.054	84	86370	6.078	ug/L	94
17) Methyl tert-butyl Ether	3.372	73	136282	5.510	ug/L	99
18) trans-1,2-Dichloroethene	3.359	96	54718	5.491	ug/L	98
19) 1,1-Dichloroethane	3.880	63	105827	5.478	ug/L	98
21) 2-Butanone	4.719	43	185062m	56.183	ug/L	
22) cis-1,2-Dichloroethene	4.674	96	59598	5.594	ug/L	98
23) Bromochloromethane	4.983	128	26285	5.596	ug/L	95
25) Chloroform	5.096	83	106664	5.382	ug/L	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU082721\
 Data File : VU044532.D
 Acq On : 27 Aug 2021 15:43
 Operator : SY/MD
 Sample : VSTDCCC005EC
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 VSTD005097

Manual Integrations
 APPROVED

MMDadoda
 8/30/2021 10:54:22 AM

Quant Time: Aug 28 01:21:21 2021
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR081821WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Sat Aug 28 01:19:29 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
27) 1,2-Dichloroethane	5.803	62	71151	5.382	ug/L	98
29) 1,1,1-Trichloroethane	5.324	97	88881	5.615	ug/L	99
30) Cyclohexane	5.398	56	92787	5.302	ug/L	99
31) Carbon tetrachloride	5.533	117	71864	5.525	ug/L	100
33) Benzene	5.784	78	237232	5.528	ug/L	100
34) Trichloroethene	6.549	95	58329	5.498	ug/L	98
35) Methylcyclohexane	6.771	83	90797	5.131	ug/L	100
37) 1,2-Dichloropropane	6.800	63	61865	5.506	ug/L	99
38) Bromodichloromethane	7.112	83	77014	5.582	ug/L	99
39) cis-1,3-Dichloropropene	7.613	75	86604	5.496	ug/L	100
40) 4-Methyl-2-pentanone	7.803	43	445064	58.334	ug/L	99
42) Toluene	7.976	91	248793	5.566	ug/L	100
44) trans-1,3-Dichloropropene	8.218	75	75269	5.635	ug/L	98
45) 1,1,2-Trichloroethane	8.407	97	46467	5.799	ug/L	97
47) Tetrachloroethene	8.562	164	39140	5.335	ug/L	98
48) 2-Hexanone	8.697	43	321914	59.132	ug/L	99
49) Dibromochloromethane	8.819	129	48608	5.732	ug/L	99
50) 1,2-Dibromoethane	8.931	107	42984	5.608	ug/L	98
51) Chlorobenzene	9.452	112	146623	5.345	ug/L	98
52) Ethylbenzene	9.578	91	259701	5.500	ug/L	99
53) m,p-Xylene	9.700	106	99844	5.624	ug/L	99
54) o-Xylene	10.108	106	97384	5.632	ug/L	99
55) Styrene	10.121	104	169399	5.975	ug/L	99
57) 1,1,2,2-Tetrachloroethane	10.790	83	59535	5.643	ug/L	98
59) Bromoform	10.298	173	25294	5.567	ug/L	98
60) Isopropylbenzene	10.491	105	254878	5.431	ug/L	99
61) 1,2,3-Trichloropropane	10.832	75	44115	5.490	ug/L	98
62) 1,3,5-Trimethylbenzene	11.095	105	208813	5.441	ug/L	99
63) 1,2,4-Trimethylbenzene	11.475	105	214631	5.454	ug/L	99
64) 1,3-Dichlorobenzene	11.751	146	110705	5.372	ug/L	100
65) 1,4-Dichlorobenzene	11.844	146	110551	5.310	ug/L	98
67) 1,2-Dichlorobenzene	12.217	146	106991	5.447	ug/L	99
68) 1,2-Dibromo-3-chloropr...	13.005	75	9165	5.576	ug/L	96
69) 1,3,5-Trichlorobenzene	13.227	180	75782	5.223	ug/L	99
70) 1,2,4-trichlorobenzene	13.848	180	59869	5.265	ug/L	99
71) Naphthalene	14.092	128	118339	5.458	ug/L	100
72) 1,2,3-Trichlorobenzene	14.336	180	57482	5.521	ug/L	99

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

