

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU020525\
 Data File : VU063165.D
 Acq On : 05 Feb 2025 17:31
 Operator : MD/SY
 Sample : Q1226-18MS
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DCZN8MS

Quant Time: Feb 06 07:31:06 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR020425WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Feb 06 07:25:52 2025
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Difluorobenzene	6.238	114	83244	5.000	ug/L	0.00
28) Chlorobenzene-d5	9.405	117	79805	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.801	152	38618	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.592	65	20004	4.201	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	84.000%
7) Chloroethane-d5	1.901	69	20878	4.724	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	94.400%
11) 1,1-Dichloroethene-d2	2.554	65	9858	4.638	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	92.800%
20) 2-Butanone-d5	4.605	46	83228	53.095	ug/L	0.00
Spiked Amount	50.000	Range	40 - 130	Recovery	=	106.200%
24) Chloroform-d	5.049	84	53077	5.042	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	100.800%
26) 1,2-Dichloroethane-d4	5.689	65	27758	5.127	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	102.600%
32) Benzene-d6	5.714	84	99380	4.774	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	95.400%
36) 1,2-Dichloropropane-d6	6.676	67	33413	4.869	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	97.400%
41) Toluene-d8	7.888	98	89612	4.873	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	97.400%
43) trans-1,3-Dichloroprop...	8.171	79	12349	4.802	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	96.000%
46) 2-Hexanone-d5	8.618	63	67465	52.128	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	104.260%
56) 1,1,2,2-Tetrachloroeth...	10.746	84	30537	5.041	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	100.800%
66) 1,2-Dichlorobenzene-d4	12.184	152	31451	4.877	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	97.600%
Target Compounds						
					Qvalue	
2) Dichlorodifluoromethane	1.377	85	37450	5.193	ug/L	99
3) Chloromethane	1.515	50	39315	4.976	ug/L	99
5) Vinyl chloride	1.599	62	41582	5.199	ug/L	99
6) Bromomethane	1.843	94	13439	3.342	ug/L	96
8) Chloroethane	1.920	64	26213	5.441	ug/L	100
9) Trichlorofluoromethane	2.126	101	49065	5.462	ug/L	99
10) 1,1,2-Trichloro-1,2,2-...	2.567	101	29502	5.343	ug/L	99
12) 1,1-Dichloroethene	2.567	96	28609	5.287	ug/L	97
13) Acetone	2.612	43	54454	55.165	ug/L	100
14) Carbon disulfide	2.779	76	98936	5.208	ug/L	99
15) Methyl Acetate	2.939	43	13433	4.983	ug/L	97
16) Methylene chloride	3.030	84	33730	5.211	ug/L	99
17) Methyl tert-butyl Ether	3.348	73	76569	5.135	ug/L	99
18) trans-1,2-Dichloroethene	3.338	96	31943	5.521	ug/L	99
19) 1,1-Dichloroethane	3.853	63	61715	5.306	ug/L	98
21) 2-Butanone	4.685	43	89366	54.527	ug/L	98
22) cis-1,2-Dichloroethene	4.650	96	81068	12.560	ug/L	99
23) Bromochloromethane	4.962	128	15678	5.415	ug/L	94
25) Chloroform	5.071	83	62213	5.434	ug/L	99

Data Path : Z:\voasrv\HPCHEM1\MSVOA_U\Data\VU020525\
 Data File : VU063165.D
 Acq On : 05 Feb 2025 17:31
 Operator : MD/SY
 Sample : Q1226-18MS
 Misc : 25.0mL/MSVOA_U/WATER
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 MSVOA_U
 ClientSampleId :
 DCZN8MS

Quant Time: Feb 06 07:31:06 2025
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_U\Method\SFAMUTR020425WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Thu Feb 06 07:25:52 2025
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
27) 1,2-Dichloroethane	5.779	62	39861	5.285	ug/L	100
29) 1,1,1-Trichloroethane	5.300	97	50410	5.337	ug/L	99
30) Cyclohexane	5.373	56	51574	5.158	ug/L	99
31) Carbon tetrachloride	5.512	117	42670	5.353	ug/L	99
33) Benzene	5.759	78	141769	5.398	ug/L	100
34) Trichloroethene	6.528	95	106652	15.990	ug/L	98
35) Methylcyclohexane	6.750	83	50957	5.098	ug/L	99
37) 1,2-Dichloropropane	6.779	63	38066	5.435	ug/L	100
38) Bromodichloromethane	7.094	83	43874	5.216	ug/L	97
39) cis-1,3-Dichloropropene	7.595	75	51379	5.144	ug/L	99
40) 4-Methyl-2-pentanone	7.775	43	219270	53.037	ug/L	99
42) Toluene	7.959	91	146556	5.339	ug/L	99
44) trans-1,3-Dichloropropene	8.200	75	43297	5.178	ug/L	100
45) 1,1,2-Trichloroethane	8.386	97	26727	5.167	ug/L	99
47) Tetrachloroethene	8.544	164	375735	80.209	ug/L	100
48) 2-Hexanone	8.669	43	159937	54.835	ug/L	99
49) Dibromochloromethane	8.798	129	29025	5.331	ug/L	98
50) 1,2-Dibromoethane	8.914	107	24960	5.261	ug/L	99
51) Chlorobenzene	9.434	112	92409	5.577	ug/L	98
52) Ethylbenzene	9.560	91	151752	5.285	ug/L	98
53) m,p-Xylene	9.682	106	56886	5.493	ug/L	99
54) o-Xylene	10.090	106	60076	5.850	ug/L	97
55) Styrene	10.103	104	90605	5.360	ug/L	96
57) 1,1,2,2-Tetrachloroethane	10.772	83	34502	5.196	ug/L	99
59) Bromoform	10.280	173	17220	5.172	ug/L	99
60) Isopropylbenzene	10.473	105	155224	5.660	ug/L	99
61) 1,2,3-Trichloropropane	10.811	75	23839	5.115	ug/L	99
62) 1,3,5-Trimethylbenzene	11.077	105	112293	5.191	ug/L	100
63) 1,2,4-Trimethylbenzene	11.457	105	134536	6.468	ug/L	99
64) 1,3-Dichlorobenzene	11.737	146	64847	5.219	ug/L	98
65) 1,4-Dichlorobenzene	11.827	146	65749	5.263	ug/L	98
67) 1,2-Dichlorobenzene	12.203	146	67850	5.844	ug/L	98
68) 1,2-Dibromo-3-chloropr...	12.984	75	5125	5.497	ug/L	98
69) 1,3,5-Trichlorobenzene	13.209	180	43120	5.079	ug/L	100
70) 1,2,4-trichlorobenzene	13.830	180	33370	4.996	ug/L	99
71) Naphthalene	14.081	128	55255	5.041	ug/L	98
72) 1,2,3-Trichlorobenzene	14.318	180	30889	5.176	ug/L	98

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

