

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW102423\
 Data File : VW032617.D
 Acq On : 25 Oct 2023 02:36
 Operator : SY/MD
 Sample : 04995-04MSD
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 A4937MSD

Manual Integrations
 APPROVED

Reviewed By : John Carlone 10/25/2023
 Supervised By : Mahesh Dadoda 10/26/2023

Quant Time: Oct 25 08:37:37 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Wed Oct 25 08:15:09 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Difluorobenzene	5.539	114	93079	5.000	ug/L	0.00
28) Chlorobenzene-d5	8.789	117	106834	5.000	ug/L	0.00
58) 1,4-Dichlorobenzene-d4	11.188	152	55369	5.000	ug/L	0.00
System Monitoring Compounds						
4) Vinyl Chloride-d3	1.278	65	23671	3.127	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	62.600%
7) Chloroethane-d5	1.532	69	27316	4.445	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	89.000%
11) 1,1-Dichloroethene-d2	2.060	65	9634	3.018	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	60.400%
20) 2-Butanone-d5	3.799	46	103690	54.147	ug/L	-0.02
Spiked Amount	50.000	Range	40 - 130	Recovery	=	108.300%
24) Chloroform-d	4.253	84	59782m	4.477	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	89.600%
26) 1,2-Dichloroethane-d4	4.947	65	31542	4.986	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	99.800%
32) Benzene-d6	4.963	84	101378	3.225	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	64.400%#
36) 1,2-Dichloropropane-d6	5.992	67	34292	3.292	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	65.800%
41) Toluene-d8	7.246	98	97524	3.393	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	67.800%#
43) trans-1,3-Dichloroprop...	7.558	79	13884	3.762	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	75.200%
46) 2-Hexanone-d5	8.027	63	90340	52.024	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	104.040%
56) 1,1,2,2-Tetrachloroeth...	10.156	84	37444	5.140	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	102.800%
66) 1,2-Dichlorobenzene-d4	11.564	152	41551	4.044	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	80.800%
Target Compounds						Qvalue
2) Dichlorodifluoromethane	1.105	85	49570	4.779	ug/L	98
3) Chloromethane	1.217	50	49330	4.718	ug/L	100
5) Vinyl chloride	1.282	62	48191	4.832	ug/L	100
6) Bromomethane	1.487	94	29373	4.876	ug/L	96
8) Chloroethane	1.548	64	34711	5.781	ug/L	97
9) Trichlorofluoromethane	1.712	101	67374	5.633	ug/L	99
10) 1,1,2-Trichloro-1,2,2-...	2.069	101	37047	5.112	ug/L	98
12) 1,1-Dichloroethene	2.069	96	33163	4.858	ug/L	88
13) Acetone	2.130	43	63553	43.059	ug/L	95
14) Carbon disulfide	2.243	76	96897	4.060	ug/L	99
15) Methyl Acetate	2.384	43	13051	3.928	ug/L	98
16) Methylene chloride	2.449	84	40234	4.521	ug/L	96
17) Methyl tert-butyl Ether	2.706	73	84814	4.564	ug/L	98
18) trans-1,2-Dichloroethene	2.696	96	33671	4.306	ug/L	98
19) 1,1-Dichloroethane	3.114	63	67713	4.604	ug/L	98
21) 2-Butanone	3.880	43	92503	40.078	ug/L	94
22) cis-1,2-Dichloroethene	3.815	96	39210	4.657	ug/L	95
23) Bromochloromethane	4.150	128	17966	5.072	ug/L	89
25) Chloroform	4.278	83	72650	5.034	ug/L	98

Data Path : Z:\voasrv\HPCHEM1\MSVOA_V\Data\VW102423\
 Data File : VW032617.D
 Acq On : 25 Oct 2023 02:36
 Operator : SY/MD
 Sample : 04995-04MSD
 Misc : 25.0mL/MSVOA_V/WATER
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 A4937MSD

Quant Time: Oct 25 08:37:37 2023
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_V\Method\SFAMVTR092923WMA.M
 Quant Title : TRACE VOA SFAM1.0
 QLast Update : Wed Oct 25 08:15:09 2023
 Response via : Initial Calibration

Manual Integrations
 APPROVED

Reviewed By :John Carlone 10/25/2023
 Supervised By :Mahesh Dadoda 10/26/2023

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
27) 1,2-Dichloroethane	5.047	62	45941	5.388	ug/L	98
29) 1,1,1-Trichloroethane	4.513	97	61691	4.247	ug/L	98
30) Cyclohexane	4.587	56	50928	3.191	ug/L	96
31) Carbon tetrachloride	4.738	117	50523	4.111	ug/L	100
33) Benzene	5.011	78	150346	3.997	ug/L	100
34) Trichloroethene	5.838	95	38253	3.725	ug/L	97
35) Methylcyclohexane	6.053	83	57784	3.636	ug/L	97
37) 1,2-Dichloropropane	6.095	63	41512	4.127	ug/L #	97
38) Bromodichloromethane	6.436	83	50706	4.242	ug/L	99
39) cis-1,3-Dichloropropene	6.957	75	59958	4.188	ug/L	99
40) 4-Methyl-2-pentanone	7.159	43	260263	40.885	ug/L	98
42) Toluene	7.317	91	182203	4.544	ug/L	99
44) trans-1,3-Dichloropropene	7.587	75	48259	4.102	ug/L	98
45) 1,1,2-Trichloroethane	7.770	97	31756	4.475	ug/L	97
47) Tetrachloroethene	7.908	164	33101	4.251	ug/L	98
48) 2-Hexanone	8.079	43	188105	41.882	ug/L	98
49) Dibromochloromethane	8.178	129	35224	4.768	ug/L	94
50) 1,2-Dibromoethane	8.284	107	29559	4.404	ug/L	98
51) Chlorobenzene	8.818	112	114474	4.434	ug/L	98
52) Ethylbenzene	8.950	91	191961	4.299	ug/L	99
53) m,p-Xylene	9.075	106	72745	4.393	ug/L	98
54) o-Xylene	9.481	106	69338	4.345	ug/L	99
55) Styrene	9.500	104	122097	4.531	ug/L	97
57) 1,1,2,2-Tetrachloroethane	10.182	83	39140	5.013	ug/L	100
59) Bromoform	9.667	173	18981	4.714	ug/L	99
60) Isopropylbenzene	9.870	105	190925	4.437	ug/L	99
61) 1,2,3-Trichloropropane	10.214	75	26677	4.527	ug/L	98
62) 1,3,5-Trimethylbenzene	10.481	105	146636	4.181	ug/L	97
63) 1,2,4-Trimethylbenzene	10.854	105	149264	4.275	ug/L	100
64) 1,3-Dichlorobenzene	11.120	146	92572	4.503	ug/L	99
65) 1,4-Dichlorobenzene	11.214	146	94760	4.595	ug/L	99
67) 1,2-Dichlorobenzene	11.580	146	86142	4.595	ug/L	99
68) 1,2-Dibromo-3-chloropr...	12.368	75	5313	4.412	ug/L	97
69) 1,3,5-Trichlorobenzene	12.583	180	65243	4.334	ug/L	99
70) 1,2,4-trichlorobenzene	13.201	180	53378	3.997	ug/L	99
71) Naphthalene	13.442	128	91449	4.031	ug/L	98
72) 1,2,3-Trichlorobenzene	13.683	180	50410	4.381	ug/L	97

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

