

Data Path : Z:\VOASRV\HPCHEM1\MSVOA V\DATA\VV071520\
 Data File : VV017501.D
 Acq On : 15 Jul 2020 10:43
 Operator : SY/MD
 Sample : L3260-01DL 4X
 Misc : 25.0mL/MSVOA V/WATER
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 MSVOA_V
 ClientSampleId :
 X2M71DL

Manual Integrations
 APPROVED

MMDadoda
 7/16/2020 4:23:54 PM

Quant Time: Jul 16 02:05:24 2020
 Quant Method : Z:\VOASRV\HPCHEM1\MSVOA_V\METHOD\SOMVTR071420WMA.M
 Quant Title : TRACE VOA SOM01.0
 QLast Update : Thu Jul 16 01:29:31 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Difluorobenzene	5.64	114	143393	5.00	ug/L	0.00
28) Chlorobenzene-d5	8.87	117	134627	5.00	ug/L	0.00
61) 1,4-Dichlorobenzene-d4	11.27	152	68922	5.00	ug/L	0.00

System Monitoring Compounds

4) Vinyl Chloride-d3	1.31	65	38528	4.62	ug/L	0.00
Spiked Amount	5.000	Range	40 - 130	Recovery	=	92.40%
7) Chloroethane-d5	1.58	69	31767	4.79	ug/L	0.00
Spiked Amount	5.000	Range	65 - 130	Recovery	=	95.80%
11) 1,1-Dichloroethene-d2	2.12	63	63889	3.50	ug/L	0.00
Spiked Amount	5.000	Range	60 - 125	Recovery	=	70.00%
20) 2-Butanone-d5	3.97	46	74456m	43.50	ug/L	0.02
Spiked Amount	50.000	Range	40 - 130	Recovery	=	87.00%
24) Chloroform-d	4.37	84	84962	4.54	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	90.80%
26) 1,2-Dichloroethane-d4	5.06	65	50307	5.06	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	101.20%
32) Benzene-d6	5.07	84	154073	4.94	ug/L	0.00
Spiked Amount	5.000	Range	70 - 125	Recovery	=	98.80%
36) 1,2-Dichloropropane-d6	6.09	67	43200	5.01	ug/L	0.00
Spiked Amount	5.000	Range	60 - 140	Recovery	=	100.20%
41) Toluene-d8	7.34	98	145142	4.78	ug/L	0.00
Spiked Amount	5.000	Range	70 - 130	Recovery	=	95.60%
45) trans-1,3-Dichloropropene-	7.65	79	19603	4.68	ug/L	0.00
Spiked Amount	5.000	Range	55 - 130	Recovery	=	93.60%
48) 2-Hexanone-d5	8.12	63	60466	48.37	ug/L	0.00
Spiked Amount	50.000	Range	45 - 130	Recovery	=	96.74%
59) 1,1,2,2-Tetrachloroethane-	10.24	84	29586	4.84	ug/L	0.00
Spiked Amount	5.000	Range	65 - 120	Recovery	=	96.80%
65) 1,2-Dichlorobenzene-d4	11.65	152	58519	4.97	ug/L	0.00
Spiked Amount	5.000	Range	80 - 120	Recovery	=	99.40%

Target Compounds

Target Compounds	R.T.	QIon	Response	Conc	Units	Ovalue
6) Bromomethane	1.53	94	3054	0.437	ug/L	92
12) 1,1-Dichloroethene	2.13	96	5098	0.591	ug/L #	1
13) Acetone	2.21	43	19677m	15.836	ug/L	
14) Carbon disulfide	2.31	76	48179	1.804	ug/L	99
15) Methyl Acetate	2.46	43	23314	9.714	ug/L	98
16) Methylene chloride	2.52	84	12287	1.183	ug/L	96
17) Methyl tert-butyl Ether	2.79	73	56356	2.629	ug/L	99
21) 2-Butanone	4.05	43	17513m	9.118	ug/L	
22) cis-1,2-Dichloroethene	3.94	96	15984	1.594	ug/L	98
27) 1,2-Dichloroethane	5.16	62	17042	1.351	ug/L #	94
30) Cyclohexane	4.70	56	35787	2.470	ug/L	99
33) Benzene	5.12	78	46965	1.311	ug/L	100
34) Trichloroethene	5.94	95	9244	0.845	ug/L	90
35) Methylcyclohexane	6.15	83	77197	4.796	ug/L	97
40) 4-Methyl-2-pentanone	7.26	43	30658	6.995	ug/L	96
42) Toluene	7.41	91	156199	3.949	ug/L	99
49) Tetrachloroethene	8.00	164	11308	1.215	ug/L	89

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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
50) 2-Hexanone	8.17	43	23884	7.417	ug/L	95
52) 1,2-Dibromoethane	8.38	107	2933	0.483	ug/L #	93
54) Ethylbenzene	9.03	91	108014	2.386	ug/L	99
55) m,p-xylene	9.16	106	34914	2.040	ug/L	100
63) 1,3-Dichlorobenzene	11.20	146	34096	1.591	ug/L	96

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

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