Data Path : Z:\SVOASRV\HPCHEM1\BNA_M\DATA\BM011720\

Quantitation Report (QT Reviewed)

Data File : BM024538.D
Acq On : 18 Jan 2020 16:35
Operator : JU
Sample : L1109-10 5X
Misc :
ALS Vial : 42 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId :

Quant Time: Jan 20 04:55:02 2020
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SOM-EPA-BM011520MA.M
Quant Title : SVOA CALIBRATION
QLast Update : Mon Jan 20 03:25:57 2020
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.46	152	132260	20.00	ng/ul	0.00
18) Naphthalene-d8	10.22	136	574751	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.10	164	549195	20.00	ng/ul	0.00
62) Phenanthrene-d10	16.87	188	1233639	20.00	ng/ul	0.02
78) Chrysene-d12	21.11	240	1879772	20.00	ng/ul	0.05
86) Perylene-d12	23.23	264	1711660	20.00	ng/ul	0.05
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0d	0.00	ng/uL	
5) Phenol-d5	6.65	99	31990	3.26	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	6.80	67	21587	3.96	ng/ul	0.00
9) 2-Chlorophenol-d4	6.99	132	27403	3.27	ng/ul	0.00
13) 4-Methylphenol-d8	8.16	113	32810	3.63	ng/ul	0.00
19) Nitrobenzene-d5	8.59	128	17057	4.22	ng/ul	0.00
22) 2-Nitrophenol-d4	9.31	143	17927	4.10	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	9.85	165	33831	3.44	ng/ul	0.00
29) 4-Chloroaniline-d4	10.35	131	236058	21.57	ng/ul	0.00
44) Dimethylphthalate-d6	13.53	166	149687	3.48	ng/ul	0.00
47) Acenaphthylene-d8	13.80	160	149521	3.02	ng/ul	0.01
52) 4-Nitrophenol-d4	14.33	143	48991	6.82	ng/ul	0.02
58) Fluorene-d10	15.12	176	136337	3.60	ng/ul	0.01
63) 4,6-Dinitro-2-methylphenol	15.23	200	10331m	1.47	ng/ul	0.00
71) Anthracene-d10	16.99	188	218411	3.81	ng/ul	0.04
79) Pyrene-d10	19.29	212	368221	3.66	ng/ul	0.03
90) Benzo(a)pyrene-d12	23.10	264	283200	3.19	ng/ul	0.06
Target Compounds						Qvalue
14) Acetophenone	8.26	105	228603m	16.112	ng/ul	
24) 2,4-Dimethylphenol	9.45	107	36401m	3.507	ng/ul	
28) Naphthalene	10.27	128	2416849	82.081	ng/ul	100
34) 2-Methylnaphthalene	11.90	142	9767251	427.492	ng/ul	99
41) 1,1'-Biphenyl	12.94	154	6682984	167.610	ng/ul	100
48) Acenaphthylene	13.83	152	15490428	306.196	ng/ul#	89
50) Acenaphthene	14.17	153	3571551	104.308	ng/ul	98
54) Dibenzofuran	14.52	168	16404284	327.037	ng/ul#	81
55) 2,4-Dinitrotoluene	14.49	165	82532	6.706	ng/ul#	66
59) Fluorene	15.18	166	18052701	425.669	ng/ul#	98
65) N-Nitrosodiphenylamine	15.40	169	358098	10.451	ng/ul#	73
70) Phenanthrene	16.91	178	37050884m	556.906	ng/ul	
72) Anthracene	17.02	178	20337811m	299.775	ng/ul	
75) Carbazole	17.28	167	10497228	183.888	ng/ul	99
77) Fluoranthene	18.95	202	30749339m	378.164	ng/ul	
80) Pyrene	19.31	202	31275893m	243.044	ng/ul	
83) Benzo(a)anthracene	21.07	228	27417434m	214.112	ng/ul	
85) Chrysene	21.13	228	22509242m	182.423	ng/ul	
88) Benzo(b)fluoranthene	22.62	252	30652314m	270.943	ng/ul	
89) Benzo(k)fluoranthene	22.65	252	8715726m	78.390	ng/ul	
91) Benzo(a)pyrene	23.06	252	15120020	153.792	ng/ul#	95
92) Indeno(1,2,3-cd)pyrene	25.35	276	11332981	110.168	ng/ul	98
93) Dibenzo(a,h)anthracene	25.34	278	4507639	51.631	ng/ul#	94

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Quant Title : SVOA CALIBRATION
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Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
94) Benzo(g,h,i)perylene	25.99	276	8719258	106.650	ng/ul	98

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

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