Quantitation Report (LSC Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM011316\
Data File : BM003978.D
Acq On : 13 Jan 2016 05:41
Operator : SJ/UM
Sample : H1095-13 10X
Misc :
ALS Vial : 20 Sample Multiplier: 1

Instrument : BNA_M
ClientSampleId : BC975

Manual Integrations APPROVED
MMdadoda
1/13/2016 3:28:50 PM

Quant Time: Jan 13 07:18:58 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-BM010116.M
Quant Title : SVOA CALIBRATION
QLast Update : Wed Jan 13 06:45:02 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	33141	20.00	ng/ul	0.00
18) Naphthalene-d8	10.45	136	165090	20.00	ng/ul	0.00
36) Acenaphthene-d10	14.32	164	102009	20.00	ng/ul	0.00
62) Phenanthrene-d10	17.07	188	235391	20.00	ng/ul	0.00
78) Chrysene-d12	21.27	240	250511	20.00	ng/ul	0.00
86) Perylene-d12	23.51	264	200424	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	0.00	96	0	0.00	ng/uL	
5) Phenol-d5	6.85	99	4980	1.81	ng/ul	0.00
7) Bis-(2-Chloroethyl)ether-d	7.00	67	3189	2.05	ng/ul	0.00
9) 2-Chlorophenol-d4	7.20	132	4500	1.99	ng/ul	0.00
13) 4-Methylphenol-d8	8.38	113	3323	1.49	ng/ul	0.00
19) Nitrobenzene-d5	8.83	128	1923	1.55	ng/ul	0.00
22) 2-Nitrophenol-d4	9.55	143	1741	1.28	ng/ul	0.00
26) 2,4-Dichlorophenol-d3	10.09	165	3779	1.56	ng/ul	0.01
29) 4-Chloroaniline-d4	10.60	131	4066	1.27	ng/ul	0.00
44) Dimethylphthalate-d6	13.73	166	16546	2.07	ng/ul	0.00
47) Acenaphthylene-d8	14.01	160	19951	2.04	ng/ul	0.00
52) 4-Nitrophenol-d4	0.00	143	0	0.00	ng/ul	
58) Fluorene-d10	15.32	176	16890	2.49	ng/ul	0.00
63) 4,6-Dinitro-2-methylphenol	15.45	200	483	0.36	ng/ul	0.00
71) Anthracene-d10	17.17	188	26421	2.43	ng/ul	0.00
79) Pyrene-d10	19.47	212	30351	2.34	ng/ul	0.00
90) Benzo(a)pyrene-d12	23.37	264	25383	2.54	ng/ul	0.00
Target Compounds						Qvalue
70) Phenanthrene	17.11	178	35672	2.69	ng/ul	99
77) Fluoranthene	19.14	202	82851	6.02	ng/ul	99
80) Pyrene	19.50	202	95250	5.60	ng/ul	98
83) Benzo(a)anthracene	21.26	228	42952	2.88	ng/ul	96
85) Chrysene	21.31	228	50665	3.57	ng/ul	98
88) Benzo(b)fluoranthene	22.84	252	50376	4.13	ng/ul#	96
89) Benzo(k)fluoranthene	22.88	252	17222m	1.49	ng/ul	
91) Benzo(a)pyrene	23.41	252	36113	3.13	ng/ul	95
92) Indeno(1,2,3-cd)pyrene	25.75	276	19502	1.53	ng/ul	99
94) Benzo(g,h,i)perylene	26.44	276	19518	1.82	ng/ul	96

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

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