Quantitation Report (LSC Reviewed)

Data Path : Z:\HPCHEM1\BNA M\DATA\BM121816\
Data File : BM008437.D
Acq On : 18 Dec 2016 14:35
Operator : UM/SJ
Sample : H6067-10DL 5X
Misc :
ALS Vial : 32 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
DA7G1DL

Manual Integrations APPROVED
umangi
12/20/2016 10:53:27 AM

Quant Time: Dec 19 02:17:57 2016
Quant Method : Z:\HPCHEM1\BNA M\METHODS\SOM02.2-EPA-2016\SOM02.2-EPA-BM121
Quant Title : SVOA CALIBRATION
QLast Update : Mon Dec 19 01:22:39 2016
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.66	152	122399	20.00	ng/ul	0.00
18) Naphthalene-d8	10.43	136	485952	20.00	ng/ul	0.00
35) Acenaphthene-d10	14.30	164	323152	20.00	ng/ul	0.00
61) Phenanthrene-d10	17.06	188	643454	20.00	ng/ul	0.00
75) Chrysene-d12	21.25	240	811080	20.00	ng/ul	0.00
83) Perylene-d12	23.48	264	856274	20.00	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.24	96	2773	1.07	ng/uL	0.01
5) Phenol-d5	6.87	99	40111	4.22	ng/ul	0.03
7) Bis-(2-Chloroethyl)ether-d	7.02	67	27489	5.04	ng/ul	0.01
9) 2-Chlorophenol-d4	7.21	132	35409	4.96	ng/ul	0.01
13) 4-Methylphenol-d8	8.40	113	28600	3.89	ng/ul	0.03
19) Nitrobenzene-d5	8.84	128	18382m	5.23	ng/ul	0.02
22) 2-Nitrophenol-d4	9.56	143	19691m	5.06	ng/ul	0.03
26) 2,4-Dichlorophenol-d3	10.13	165	34546m	4.84	ng/ul	0.06
29) 4-Chloroaniline-d4	10.65	131	23587m	2.89	ng/ul	0.06
43) Dimethylphthalate-d6	13.73	166	132259	6.16	ng/ul	0.01
46) Acenaphthylene-d8	13.99	160	157506	6.20	ng/ul	0.00
51) 4-Nitrophenol-d4	14.70	143	2416m	0.76	ng/ul	0.15
57) Fluorene-d10	15.31	176	114511	6.13	ng/ul	0.01
62) 4,6-Dinitro-2-methylphenol	15.49	200	7721	2.02	ng/ul	0.04
70) Anthracene-d10	17.16	188	176943	6.31	ng/ul	0.00
76) Pyrene-d10	19.46	212	213502	6.82	ng/ul	0.00
87) Benzo(a)pyrene-d12	23.34	264	224634	6.25	ng/ul	0.00
Target Compounds						Qvalue
6) Phenol	6.90	94	9980	1.01	ng/ul#	83
44) Dimethylphthalate	13.77	163	34460	1.64	ng/ul	96
47) Acenaphthylene	14.03	152	33935	1.31	ng/ul	99
71) Anthracene	17.20	178	42665	1.27	ng/ul	96
77) Pyrene	19.49	202	49136	1.23	ng/ul	98
80) Benzo(a)anthracene	21.24	228	76985m	1.84	ng/ul	
81) Bis(2-ethylhexyl)phthalate	21.16	149	692184	33.09	ng/ul	99
82) Chrysene	21.29	228	157802	3.91	ng/ul	95
85) Benzo(b)fluoranthene	22.81	252	452938	9.80	ng/ul	97
86) Benzo(k)fluoranthene	22.86	252	132710m	3.00	ng/ul	
88) Benzo(a)pyrene	23.39	252	311485	6.99	ng/ul	96
89) Indeno(1,2,3-cd)pyrene	25.71	276	324926	6.01	ng/ul	97
90) Dibenzo(a,h)anthracene	25.71	278	86750	1.91	ng/ul	92
91) Benzo(g,h,i)perylene	26.40	276	293535	6.49	ng/ul	100

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

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