Quantitation Report (Not Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE063015\
Data File : BE090038.D
Acq On : 1 Jul 2015 8:58
Operator : TP/IZ
Sample : G2824-01 5X
Misc :
ALS Vial : 24 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : X1SS0650006-150626

Quant Time: Jul 01 17:23:13 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062915.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Jun 30 07:01:01 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	9	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	6	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	37	5.00	ng	-0.02
25) Chrysene-d12	21.13	240	78	5.00	ng	0.00
32) Perylene-d12	23.44	264	198	5.00	ng	-0.01
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	1	0.87	ng	0.01
5) Phenol-d6	6.79	99	1	0.62	ng	0.00
7) Nitrobenzene-d5	8.74	82	3	4.20	ng	-0.02
13) 2,4,6-Tribromophenol	15.76	330	1	5.69	ng	0.03
14) 2-Fluorobiphenyl	12.83	172	2	1.11	ng	-0.02
27) Terphenyl-d14	19.58	244	22	1.69	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	6	8.74	ng	# 19
3) n-Nitrosodimethylamine	3.55	42	6	6.45	ng	# 1
8) Nitrobenzene	8.78	77	4	5.27	ng	# 63
9) Naphthalene	10.42	128	5	2.46	ng	# 1
10) Hexachlorobutadiene	10.55	225	4	12.51	ng	85
11) 2-Methylnaphthalene	12.03	142	4	2.94	ng	# 55
15) Acenaphthylene	13.95	152	4	0.37	ng	# 60
16) Acenaphthene	14.33	154	4	2.57	ng	# 1
17) Fluorene	15.33	166	17	8.01	ng	# 1
19) 4-Bromophenyl-phenylether	16.20	248	15	9.54	ng	# 69
20) Hexachlorobenzene	16.27	284	6	0.92	ng	# 36
21) Pentachlorophenol	16.65	266	7	8.34	ng	# 55
22) Phenanthrene	17.01	178	66	6.96	ng	# 71
23) Anthracene	17.01	178	54	6.23	ng	# 66
24) Fluoranthene	19.01	202	60	5.67	ng	# 78
26) Pyrene	19.37	202	58	3.12	ng	94
28) Benzo(a)anthracene	21.11	228	83	4.26	ng	# 8
29) Chrysene	21.15	228	85	4.20	ng	# 17
30) Bis(2-ethylhexyl)phthalate	21.04	149	9	1.25	ng	# 54
31) Indeno(1,2,3-cd)pyrene	25.83	276	297	15.31	ng	# 80
33) Benzo(b)fluoranthene	22.73	252	330	7.55	ng	# 6
34) Benzo(k)fluoranthene	22.73	252	334	6.84	ng	# 6
35) Benzo(a)pyrene	23.33	252	218	5.40	ng	# 1
36) Dibenzo(a,h)anthracene	25.86	278	121	2.92	ng	# 1
37) Benzo(g,h,i)perylene	26.58	276	309	7.36	ng	# 33

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

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