Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070615\
Data File : BE090060.D
Acq On : 6 Jul 2015 19:18
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Jul 07 01:17:15 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	7.61	152	23959	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	104080	5.00	ng	-0.01
12) Acenaphthene-d10	14.22	164	58377	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	133688	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	154158	5.00	ng	-0.01
32) Perylene-d12	23.43	264	141308	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	5065	0.92	ng	0.00
5) Phenol-d6	6.80	99	7323	0.95	ng	0.00
7) Nitrobenzene-d5	8.75	82	7097	0.86	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	1350	0.79	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	15940	0.91	ng	-0.01
27) Terphenyl-d14	19.58	244	23680	0.92	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	2676	0.81	ng	97
3) n-Nitrosodimethylamine	3.52	42	3633	0.82	ng	94
8) Nitrobenzene	8.79	77	7487	0.85	ng	99
9) Naphthalene	10.41	128	21669	0.92	ng	99
10) Hexachlorobutadiene	10.72	225	3491	0.94	ng	99
11) 2-Methylnaphthalene	12.03	142	13870	0.88	ng	100
15) Acenaphthylene	13.93	152	96637	0.92	ng	100
16) Acenaphthene	14.29	154	14145	0.93	ng	98
17) Fluorene	15.27	166	18347	0.89	ng	100
19) 4-Bromophenyl-phenylether	16.16	248	4944	0.87	ng	94
20) Hexachlorobenzene	16.28	284	21186	0.90	ng	97
21) Pentachlorophenol	16.63	266	1467	0.73	ng	96
22) Phenanthrene	17.00	178	29904	0.87	ng	100
23) Anthracene	17.10	178	28192	0.90	ng	100
24) Fluoranthene	19.00	202	33656	0.88	ng	100
26) Pyrene	19.36	202	34978	0.95	ng	100
28) Benzo(a)anthracene	21.10	228	35381	0.92	ng	99
29) Chrysene	21.15	228	36059	0.90	ng	99
30) Bis(2-ethylhexyl)phthalate	21.05	149	10620	0.87	ng	99
31) Indeno(1,2,3-cd)pyrene	25.83	276	35074	0.92	ng	99
33) Benzo(b)fluoranthene	22.73	252	31802	1.02	ng	100
34) Benzo(k)fluoranthene	22.77	252	34154	0.98	ng	99
35) Benzo(a)pyrene	23.32	252	29097	1.01	ng	99
36) Dibenzo(a,h)anthracene	25.85	278	28084	0.95	ng	98
37) Benzo(g,h,i)perylene	26.57	276	28394	0.95	ng	98

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

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