Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070115\
Data File : BE090042.D
Acq On : 1 Jul 2015 12:45
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Jul 02 06:34:21 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	20373	5.00	ng	-0.02
6) Naphthalene-d8	10.37	136	85773	5.00	ng	-0.01
12) Acenaphthene-d10	14.22	164	48832	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	114468	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	138511	5.00	ng	-0.01
32) Perylene-d12	23.44	264	129459	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	4237	0.91	ng	-0.01
5) Phenol-d6	6.79	99	6002	0.92	ng	-0.01
7) Nitrobenzene-d5	8.75	82	5820	0.85	ng	-0.01
13) 2,4,6-Tribromophenol	15.73	330	1112	0.78	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	13280	0.90	ng	-0.01
27) Terphenyl-d14	19.58	244	20528	0.89	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	2400	0.86	ng	95
3) n-Nitrosodimethylamine	3.51	42	3097	0.82	ng	94
8) Nitrobenzene	8.79	77	6199	0.86	ng	98
9) Naphthalene	10.41	128	17861	0.92	ng	99
10) Hexachlorobutadiene	10.71	225	2894	0.95	ng	99
11) 2-Methylnaphthalene	12.03	142	11574	0.89	ng	98
15) Acenaphthylene	13.94	152	81238	0.92	ng	100
16) Acenaphthene	14.29	154	11988	0.94	ng	98
17) Fluorene	15.27	166	15999	0.93	ng	99
19) 4-Bromophenyl-phenylether	16.16	248	4354	0.89	ng	98
20) Hexachlorobenzene	16.28	284	18534	0.92	ng	97
21) Pentachlorophenol	16.63	266	1240	0.72	ng	96
22) Phenanthrene	17.00	178	26040	0.89	ng	100
23) Anthracene	17.10	178	24625	0.92	ng	99
24) Fluoranthene	19.01	202	29511	0.90	ng	100
26) Pyrene	19.37	202	30649	0.93	ng	100
28) Benzo(a)anthracene	21.10	228	31594	0.91	ng	99
29) Chrysene	21.16	228	32225	0.90	ng	99
30) Bis(2-ethylhexyl)phthalate	21.05	149	9786	0.88	ng	100
31) Indeno(1,2,3-cd)pyrene	25.84	276	32741	0.95	ng	98
33) Benzo(b)fluoranthene	22.73	252	29365	1.03	ng	100
34) Benzo(k)fluoranthene	22.78	252	31555	0.99	ng	100
35) Benzo(a)pyrene	23.33	252	26691	1.01	ng	99
36) Dibenzo(a,h)anthracene	25.86	278	26309	0.97	ng	99
37) Benzo(g,h,i)perylene	26.58	276	26588	0.97	ng	98

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

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