Quantitation Report (Not Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061217\
Data File : BE093095.D
Acq On : 12 Jun 2017 9:08
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: Jun 12 19:28:33 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE061217.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Mon Jun 12 14:07:54 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.75	152	731	0.40	ng		0.00
7) Naphthalene-d8	10.52	136	3437	0.40	ng		0.00
13) Acenaphthene-d10	14.37	164	1702	0.40	ng		0.00
19) Phenanthrene-d10	17.13	188	3503	0.40	ng		0.00
27) Chrysene-d12	21.27	240	5607	0.40	ng		0.00
34) Perylene-d12	23.67	264	4950	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.34	112	769	0.36	ng		0.00
5) Phenol-d6	6.90	99	1103	0.35	ng		0.00
8) Nitrobenzene-d5	8.88	82	981	0.36	ng		0.00
11) 2-Methylnaphthalene-d10	12.11	152	1850	0.36	ng		0.00
14) 2,4,6-Tribromophenol	15.85	330	133	0.34	ng		0.00
15) 2-Fluorobiphenyl	12.99	172	2662	0.37	ng		0.00
25) Fluoranthene-d10	19.11	212	4093	0.35	ng		0.00
29) Terphenyl-d14	19.72	244	3936	0.36	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.29	88	558	0.344	ng		99
3) n-Nitrosodimethylamine	3.60	42	495	0.357	ng	#	99
6) bis(2-Chloroethyl)ether	7.15	93	985	0.357	ng	#	99
9) Naphthalene	10.56	128	3296	0.357	ng		98
10) Hexachlorobutadiene	10.87	225	437	0.349	ng		99
12) 2-Methylnaphthalene	12.18	142	2024	0.354	ng		97
16) Acenaphthylene	14.08	152	10242	0.346	ng		100
17) Acenaphthene	14.44	154	2007	0.364	ng		97
18) Fluorene	15.41	166	2397	0.360	ng		99
20) 4-Bromophenyl-phenylether	16.33	248	777	0.364	ng	#	80
21) Hexachlorobenzene	16.44	284	671	0.358	ng	#	100
22) Pentachlorophenol	16.79	266	214	0.354	ng		89
23) Phenanthrene	17.16	178	5840	0.369	ng		99
24) Anthracene	17.25	178	4469	0.333	ng		99
26) Fluoranthene	19.16	202	20543	0.357	ng	#	99
28) Pyrene	19.51	202	26404	0.387	ng	#	98
30) Benzo(a)anthracene	21.23	228	5542	0.366	ng	#	85
31) Chrysene	21.28	228	5703	0.353	ng	#	84
32) Bis(2-ethylhexyl)phthalate	21.18	149	2231	0.359	ng		96
33) Indeno(1,2,3-cd)pyrene	26.17	276	21579	0.357	ng		99
35) Benzo(b)fluoranthene	22.93	252	5417	0.341	ng		97
36) Benzo(k)fluoranthene	22.97	252	5317	0.347	ng		98
37) Benzo(a)pyrene	23.55	252	4637	0.337	ng		97
38) Dibenzo(a,h)anthracene	26.20	278	4954	0.346	ng		96
39) Benzo(g,h,i)perylene	26.95	276	19789	0.342	ng		97

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

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