Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE051017\
Data File : BE092979.D
Acq On : 10 May 2017 13:21
Operator : SJ/MA
Sample : SSTDCCC0.4
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC0.4

Quant Time: May 10 14:08:30 2017
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE051017.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed May 10 03:43:15 2017
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.77	152	795	0.40	ng		0.00
7) Naphthalene-d8	10.53	136	3351	0.40	ng		0.00
13) Acenaphthene-d10	14.39	164	1645	0.40	ng		0.00
19) Phenanthrene-d10	17.13	188	3992	0.40	ng		0.00
26) Chrysene-d12	21.28	240	4155	0.40	ng		0.00
33) Perylene-d12	23.70	264	3623	0.40	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.35	112	833	0.37	ng		0.00
5) Phenol-d6	6.95	99	1048	0.38	ng		0.00
8) Nitrobenzene-d5	8.90	82	930	0.36	ng		0.00
11) 2-Methylnaphthalene-d10	12.13	152	1804	0.38	ng		0.00
14) 2,4,6-Tribromophenol	15.88	330	138	0.32	ng		0.00
15) 2-Fluorobiphenyl	13.01	172	2444	0.40	ng		0.00
24) Fluoranthene-d10	19.15	212	3893	0.37	ng		0.00
28) Terphenyl-d14	19.74	244	2428	0.38	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.31	88	631	0.38	ng		97
3) n-Nitrosodimethylamine	3.60	42	547	0.40	ng		98
6) bis(2-Chloroethyl)ether	7.20	93	1103	0.41	ng	#	100
9) Naphthalene	10.58	128	3416	0.38	ng		98
10) Hexachlorobutadiene	10.89	225	485	0.38	ng		99
12) 2-Methylnaphthalene	12.19	142	1991	0.39	ng		91
16) Acenaphthylene	14.10	152	10066	0.36	ng		100
17) Acenaphthene	14.46	154	1970	0.38	ng		99
18) Fluorene	15.44	166	2399	0.38	ng		99
20) Hexachlorobenzene	16.47	284	731	0.39	ng	#	100
21) Pentachlorophenol	16.81	266	148	0.37	ng		98
22) Phenanthrene	17.18	178	4730	0.42	ng		99
23) Anthracene	17.27	178	3569	0.40	ng		99
25) Fluoranthene	19.17	202	20601	0.38	ng	#	99
27) Pyrene	19.53	202	22052	0.40	ng	#	98
29) Benzo(a)anthracene	21.27	228	3766	0.38	ng		99
30) Chrysene	21.32	228	4922	0.37	ng		99
31) Bis(2-ethylhexyl)phthalate	21.20	149	1574	0.41	ng		97
32) Indeno(1,2,3-cd)pyrene	26.24	276	16633	0.36	ng		100
34) Benzo(b)fluoranthene	22.95	252	3891	0.35	ng		100
35) Benzo(k)fluoranthene	23.01	252	4476	0.38	ng		99
36) Benzo(a)pyrene	23.59	252	3563	0.34	ng		99
37) Dibenzo(a,h)anthracene	26.25	278	3610	0.35	ng		99
38) Benzo(g,h,i)perylene	27.00	276	15981	0.36	ng		100

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

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