Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089985.D
Acq On : 29 Jun 2015 16:09
Operator : TP/IZ
Sample : SSTDICCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICCC001

Quant Time: Jun 30 06:06:20 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 06:02:58 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	20316	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	86161	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	51874	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	128935	5.00	ng	0.00
25) Chrysene-d12	21.13	240	159817	5.00	ng	0.00
32) Perylene-d12	23.46	264	154962	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	4654	1.22	ng	0.00
5) Phenol-d6	6.80	99	6421	1.11	ng	0.00
7) Nitrobenzene-d5	8.76	82	6722	1.03	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	1488	1.90	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	15180	0.99	ng	0.00
27) Terphenyl-d14	19.59	244	26582	0.96	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	2684	0.96	ng	100
3) n-Nitrosodimethylamine	3.53	42	3629	0.98	ng	100
8) Nitrobenzene	8.80	77	7033	1.01	ng	100
9) Naphthalene	10.43	128	19150	0.99	ng	100
10) Hexachlorobutadiene	10.72	225	3014	0.99	ng	100
11) 2-Methylnaphthalene	12.04	142	12781	0.97	ng	100
15) Acenaphthylene	13.95	152	90725	0.99	ng	100
16) Acenaphthene	14.30	154	13283	0.99	ng	100
17) Fluorene	15.29	166	17859	0.98	ng	100
19) 4-Bromophenyl-phenylether	16.18	248	5358	1.00	ng	100
20) Hexachlorobenzene	16.30	284	21942	0.95	ng	100
21) Pentachlorophenol	16.63	266	2094	1.96	ng	100
22) Phenanthrene	17.01	178	32055	0.98	ng	100
23) Anthracene	17.10	178	29649	0.98	ng	100
24) Fluoranthene	19.02	202	36164	1.01	ng	100
26) Pyrene	19.38	202	37429	0.93	ng	100
28) Benzo(a)anthracene	21.11	228	39021	1.01	ng	100
29) Chrysene	21.17	228	40942	0.97	ng	100
30) Bis(2-ethylhexyl)phthalate	21.07	149	11691	2.46	ng	100
31) Indeno(1,2,3-cd)pyrene	25.87	276	39084	1.22	ng	100
33) Benzo(b)fluoranthene	22.75	252	33945	1.02	ng	100
34) Benzo(k)fluoranthene	22.79	252	38102	1.01	ng	100
35) Benzo(a)pyrene	23.35	252	31475	1.06	ng	100
36) Dibenzo(a,h)anthracene	25.89	278	31835	1.14	ng	100
37) Benzo(g,h,i)perylene	26.61	276	32480	1.09	ng	100

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

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