Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
Data File : BE090074.D
Acq On : 9 Jul 2015 00:34
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Jul 09 01:21:53 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	29794	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	132416	5.00	ng	-0.01
12) Acenaphthene-d10	14.23	164	71854	5.00	ng	0.00
18) Phenanthrene-d10	16.96	188	159228	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	182232	5.00	ng	-0.01
32) Perylene-d12	23.43	264	170648	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	6955	1.02	ng	0.00
5) Phenol-d6	6.80	99	10312	1.08	ng	0.00
7) Nitrobenzene-d5	8.75	82	9586	0.91	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	2281	1.08	ng	0.00
14) 2-Fluorobiphenyl	12.85	172	19935	0.92	ng	0.00
27) Terphenyl-d14	19.58	244	29150	0.96	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	3112	0.76	ng	97
3) n-Nitrosodimethylamine	3.53	42	4496	0.81	ng	90
8) Nitrobenzene	8.79	77	9870	0.88	ng	97
9) Naphthalene	10.41	128	27583	0.92	ng	99
10) Hexachlorobutadiene	10.72	225	4413	0.94	ng	100
11) 2-Methylnaphthalene	12.03	142	17804	0.89	ng	99
15) Acenaphthylene	13.94	152	123143	0.95	ng	100
16) Acenaphthene	14.29	154	17624	0.94	ng	97
17) Fluorene	15.27	166	22661	0.89	ng	99
19) 4-Bromophenyl-phenylether	16.16	248	6138	0.91	ng	95
20) Hexachlorobenzene	16.28	284	25178	0.89	ng	98
21) Pentachlorophenol	16.63	266	2706	0.98	ng	99
22) Phenanthrene	17.00	178	37076	0.91	ng	100
23) Anthracene	17.10	178	34627	0.93	ng	99
24) Fluoranthene	19.00	202	41707	0.92	ng	100
26) Pyrene	19.37	202	43481	1.00	ng	100
28) Benzo(a)anthracene	21.10	228	43656	0.96	ng	99
29) Chrysene	21.15	228	43392	0.92	ng	99
30) Bis(2-ethylhexyl)phthalate	21.05	149	21932	1.29	ng	98
31) Indeno(1,2,3-cd)pyrene	25.83	276	45561	1.01	ng	98
33) Benzo(b)fluoranthene	22.73	252	40438	1.07	ng	99
34) Benzo(k)fluoranthene	22.77	252	42916	1.02	ng	99
35) Benzo(a)pyrene	23.33	252	37584	1.08	ng	99
36) Dibenzo(a,h)anthracene	25.85	278	36296	1.02	ng	98
37) Benzo(g,h,i)perylene	26.57	276	37226	1.03	ng	97

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

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