Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE070815\
Data File : BE090064.D
Acq On : 8 Jul 2015 18:30
Operator : TP/IZ
Sample : PB84399BS
Misc :
ALS Vial : 4 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : PB84399BS

Quant Time: Jul 09 01:19:31 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	7.62	152	15837	5.00	ng	0.00
6) Naphthalene-d8	10.37	136	69989	5.00	ng	-0.01
12) Acenaphthene-d10	14.22	164	39669	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	87819	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	106604	5.00	ng	-0.01
32) Perylene-d12	23.43	264	92954	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	20279	5.57	ng	0.00
5) Phenol-d6	6.80	99	30958	6.08	ng	0.00
7) Nitrobenzene-d5	8.75	82	20274	3.65	ng	-0.01
13) 2,4,6-Tribromophenol	15.73	330	7194	6.19	ng	0.00
14) 2-Fluorobiphenyl	12.84	172	47275	3.97	ng	-0.01
27) Terphenyl-d14	19.58	244	71976	4.05	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	6354	2.92	ng	97
3) n-Nitrosodimethylamine	3.52	42	8716	2.96	ng	99
8) Nitrobenzene	8.79	77	19852	3.37	ng	98
9) Naphthalene	10.42	128	52297	3.30	ng	99
10) Hexachlorobutadiene	10.72	225	8193	3.29	ng	99
11) 2-Methylnaphthalene	12.03	142	36535	3.46	ng	99
15) Acenaphthylene	13.94	152	253027	3.54	ng	100
16) Acenaphthene	14.29	154	36979	3.59	ng	97
17) Fluorene	15.27	166	49720	3.54	ng	99
19) 4-Bromophenyl-phenylether	16.16	248	14315	3.83	ng	95
20) Hexachlorobenzene	16.28	284	60118	3.87	ng	97
21) Pentachlorophenol	16.63	266	9719	4.98	ng	96
22) Phenanthrene	17.00	178	83177	3.70	ng	100
23) Anthracene	17.08	178	78595	3.82	ng	99
24) Fluoranthene	19.00	202	94937	3.78	ng	100
26) Pyrene	19.36	202	100297	3.95	ng	99
28) Benzo(a)anthracene	21.10	228	98598	3.70	ng	99
29) Chrysene	21.15	228	103536	3.75	ng	100
30) Bis(2-ethylhexyl)phthalate	21.05	149	28594	2.50	ng	99
31) Indeno(1,2,3-cd)pyrene	25.84	276	91921	3.47	ng	99
33) Benzo(b)fluoranthene	22.73	252	86884	4.24	ng	99
34) Benzo(k)fluoranthene	22.77	252	95357	4.16	ng	99
35) Benzo(a)pyrene	23.33	252	82169	4.34	ng	99
36) Dibenzo(a,h)anthracene	25.85	278	75136	3.86	ng	99
37) Benzo(g,h,i)perylene	26.57	276	76438	3.88	ng	99

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

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