Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062415\
Data File : BE089940.D
Acq On : 24 Jun 2015 19:38
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Jun 25 03:54:08 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Jun 24 09:27:42 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	27506	5.00	ng		0.00
6) Naphthalene-d8	10.39	136	127771	5.00	ng		0.00
12) Acenaphthene-d10	14.25	164	78547	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	189976	5.00	ng		0.00
25) Chrysene-d12	21.14	240	199601	5.00	ng		0.00
32) Perylene-d12	23.47	264	171293	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	3931	0.79	ng		0.00
5) Phenol-d6	6.81	99	6204	0.82	ng		0.00
7) Nitrobenzene-d5	8.77	82	8194	0.86	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	651	0.77	ng		0.00
14) 2-Fluorobiphenyl	12.87	172	20749	0.91	ng		0.00
27) Terphenyl-d14	19.60	244	31248	0.91	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	3225	0.97	ng		96
3) n-Nitrosodimethylamine	3.55	42	4402	0.89	ng		93
8) Nitrobenzene	8.81	77	8735	0.86	ng		99
9) Naphthalene	10.44	128	26214	0.92	ng		99
10) Hexachlorobutadiene	10.74	225	4217	0.94	ng		99
11) 2-Methylnaphthalene	12.05	142	17477	0.90	ng		98
15) Acenaphthylene	13.96	152	127862	0.93	ng		100
16) Acenaphthene	14.30	154	19216	0.95	ng		98
17) Fluorene	15.29	166	25275	0.99	ng		98
19) 4-Bromophenyl-phenylether	16.20	248	6853	0.88	ng		97
20) Hexachlorobenzene	16.30	284	31009	0.91	ng		99
21) Pentachlorophenol	16.63	266	805	0.78	ng	#	92
22) Phenanthrene	17.03	178	43162	0.91	ng		99
23) Anthracene	17.12	178	38931	0.88	ng		100
24) Fluoranthene	19.02	202	46278	0.89	ng		100
26) Pyrene	19.38	202	47642	0.94	ng		100
28) Benzo(a)anthracene	21.12	228	42629	0.89	ng		99
29) Chrysene	21.18	228	46693	0.90	ng		99
30) Bis(2-ethylhexyl)phthalate	21.07	149	3101	0.76	ng		98
31) Indeno(1,2,3-cd)pyrene	25.89	276	34425	0.90	ng	#	100
33) Benzo(b)fluoranthene	22.76	252	34004	0.94	ng		99
34) Benzo(k)fluoranthene	22.80	252	38496	0.94	ng		100
35) Benzo(a)pyrene	23.36	252	29085	0.90	ng		99
36) Dibenzo(a,h)anthracene	25.91	278	27147	0.91	ng		99
37) Benzo(g,h,i)perylene	26.63	276	29617	0.92	ng		99

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

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