Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062415\
Data File : BE089933.D
Acq On : 24 Jun 2015 14:30
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001

Quant Time: Jun 25 03:45:05 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	29491	5.00	ng		0.00
6) Naphthalene-d8	10.39	136	136189	5.00	ng		0.00
12) Acenaphthene-d10	14.24	164	83247	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	195463	5.00	ng		0.00
25) Chrysene-d12	21.14	240	219183	5.00	ng		0.00
32) Perylene-d12	23.47	264	192639	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	4420	0.83	ng		0.00
5) Phenol-d6	6.81	99	6938	0.85	ng		0.00
7) Nitrobenzene-d5	8.77	82	9047	0.89	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	737	0.80	ng		0.00
14) 2-Fluorobiphenyl	12.87	172	21815	0.90	ng		0.00
27) Terphenyl-d14	19.60	244	33706	0.89	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	3327	0.93	ng		99
3) n-Nitrosodimethylamine	3.55	42	4723	0.89	ng	#	92
8) Nitrobenzene	8.81	77	9553	0.89	ng		100
9) Naphthalene	10.44	128	28054	0.93	ng		99
10) Hexachlorobutadiene	10.74	225	4436	0.93	ng		99
11) 2-Methylnaphthalene	12.06	142	18479	0.89	ng		99
15) Acenaphthylene	13.96	152	134730	0.92	ng		100
16) Acenaphthene	14.30	154	20102	0.94	ng		98
17) Fluorene	15.29	166	26352	0.98	ng		99
19) 4-Bromophenyl-phenylether	16.20	248	7108	0.89	ng		98
20) Hexachlorobenzene	16.30	284	32352	0.93	ng		99
21) Pentachlorophenol	16.63	266	878	0.80	ng		90
22) Phenanthrene	17.03	178	45253	0.92	ng		100
23) Anthracene	17.12	178	41494	0.91	ng		99
24) Fluoranthene	19.02	202	49155	0.92	ng		100
26) Pyrene	19.38	202	50951	0.92	ng		100
28) Benzo(a)anthracene	21.12	228	47332	0.90	ng		99
29) Chrysene	21.17	228	50991	0.89	ng		98
30) Bis(2-ethylhexyl)phthalate	21.07	149	3506	0.79	ng		100
31) Indeno(1,2,3-cd)pyrene	25.88	276	38852	0.92	ng	#	100
33) Benzo(b)fluoranthene	22.75	252	38460	0.94	ng		99
34) Benzo(k)fluoranthene	22.80	252	42429	0.92	ng		100
35) Benzo(a)pyrene	23.36	252	32869	0.91	ng		99
36) Dibenzo(a,h)anthracene	25.91	278	30525	0.91	ng		99
37) Benzo(g,h,i)perylene	26.62	276	33274	0.92	ng		100

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

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