Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089995.D
Acq On : 29 Jun 2015 22:43
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Jun 30 07:03:48 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	7.63	152	18452	5.00	ng		0.00
6) Naphthalene-d8	10.38	136	77244	5.00	ng		0.00
12) Acenaphthene-d10	14.23	164	45593	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	112365	5.00	ng		0.00
25) Chrysene-d12	21.13	240	142268	5.00	ng		0.00
32) Perylene-d12	23.46	264	133254	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	3408	0.80	ng		0.00
5) Phenol-d6	6.80	99	4857	0.82	ng		0.00
7) Nitrobenzene-d5	8.76	82	5385	0.88	ng		0.00
13) 2,4,6-Tribromophenol	15.73	330	835	0.62	ng		0.00
14) 2-Fluorobiphenyl	12.86	172	13436	0.98	ng		0.00
27) Terphenyl-d14	19.59	244	22709	0.96	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.24	88	2565	1.01	ng		94
3) n-Nitrosodimethylamine	3.53	42	3137	0.91	ng	#	93
8) Nitrobenzene	8.80	77	5779	0.89	ng		98
9) Naphthalene	10.43	128	16789	0.96	ng		99
10) Hexachlorobutadiene	10.72	225	2785	1.01	ng		99
11) 2-Methylnaphthalene	12.04	142	11106	0.95	ng		99
15) Acenaphthylene	13.95	152	77947	0.95	ng		100
16) Acenaphthene	14.29	154	11510	0.97	ng		97
17) Fluorene	15.29	166	15368	0.95	ng		99
19) 4-Bromophenyl-phenylether	16.18	248	4656	0.97	ng		99
20) Hexachlorobenzene	16.30	284	19458	0.98	ng		99
21) Pentachlorophenol	16.63	266	1076	0.67	ng		93
22) Phenanthrene	17.01	178	28086	0.98	ng		100
23) Anthracene	17.10	178	24744	0.94	ng		99
24) Fluoranthene	19.01	202	30924	0.96	ng		100
26) Pyrene	19.38	202	32168	0.95	ng		100
28) Benzo(a)anthracene	21.11	228	33200	0.93	ng		100
29) Chrysene	21.16	228	35973	0.98	ng		98
30) Bis(2-ethylhexyl)phthalate	21.07	149	5339	0.61	ng		100
31) Indeno(1,2,3-cd)pyrene	25.87	276	30928	0.87	ng		100
33) Benzo(b)fluoranthene	22.75	252	27957	0.95	ng		99
34) Benzo(k)fluoranthene	22.79	252	32313	0.98	ng		99
35) Benzo(a)pyrene	23.35	252	25081	0.92	ng		99
36) Dibenzo(a,h)anthracene	25.89	278	25169	0.90	ng		99
37) Benzo(g,h,i)perylene	26.61	276	25464	0.90	ng		100

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

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