Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
Data File : BE089962.D
Acq On : 27 Jun 2015 4:34
Operator : TP/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 2 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Jun 27 06:30:22 2015
Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Thu Jun 25 16:51:27 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	25185	5.00	ng		0.00
6) Naphthalene-d8	10.39	136	124716	5.00	ng		0.00
12) Acenaphthene-d10	14.24	164	77301	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	179295	5.00	ng		0.00
25) Chrysene-d12	21.13	240	180415	5.00	ng		0.00
32) Perylene-d12	23.47	264	149956	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	6194	1.36	ng		0.00
5) Phenol-d6	6.81	99	16054	2.30	ng		0.00
7) Nitrobenzene-d5	8.77	82	12951	1.39	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	1997	1.72	ng		0.00
14) 2-Fluorobiphenyl	12.86	172	40445	1.80	ng		0.00
27) Terphenyl-d14	19.60	244	60285	1.94	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	3216	1.05	ng		89
3) n-Nitrosodimethylamine	3.55	42	3749	0.83	ng	#	91
8) Nitrobenzene	8.81	77	7474	0.76	ng		100
9) Naphthalene	10.44	128	23025	0.83	ng		99
10) Hexachlorobutadiene	10.74	225	3753	0.86	ng		100
11) 2-Methylnaphthalene	12.05	142	14956	0.79	ng		98
15) Acenaphthylene	13.95	152	106338	0.78	ng		99
16) Acenaphthene	14.30	154	16087	0.81	ng		97
17) Fluorene	15.29	166	20784	0.83	ng		99
19) 4-Bromophenyl-phenylether	16.18	248	5511	0.75	ng		99
20) Hexachlorobenzene	16.30	284	26335	0.82	ng		99
21) Pentachlorophenol	16.63	266	769	0.85	ng	#	90
22) Phenanthrene	17.01	178	35148	0.78	ng		100
23) Anthracene	17.12	178	31650	0.76	ng		100
24) Fluoranthene	19.02	202	36114	0.74	ng		99
26) Pyrene	19.39	202	37575	0.82	ng		100
28) Benzo(a)anthracene	21.12	228	32881	0.76	ng		98
29) Chrysene	21.17	228	35709	0.76	ng		98
30) Bis(2-ethylhexyl)phthalate	21.07	149	2781	0.76	ng		99
31) Indeno(1,2,3-cd)pyrene	25.88	276	25811	0.75	ng	#	100
33) Benzo(b)fluoranthene	22.76	252	25480	0.80	ng		100
34) Benzo(k)fluoranthene	22.80	252	29038	0.81	ng		99
35) Benzo(a)pyrene	23.36	252	21542	0.76	ng		100
36) Dibenzo(a,h)anthracene	25.90	278	20465	0.78	ng		99
37) Benzo(g,h,i)perylene	26.62	276	21871	0.78	ng		99

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

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