Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089988.D
Acq On : 29 Jun 2015 17:56
Operator : TP/IZ
Sample : SSTDICC0.2
Misc :
ALS Vial : 9 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.2

Quant Time: Jun 30 06:07:17 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 06:02:58 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	25044	5.00	ng		0.00
6) Naphthalene-d8	10.38	136	104269	5.00	ng		0.00
12) Acenaphthene-d10	14.23	164	61272	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	155015	5.00	ng		0.00
25) Chrysene-d12	21.13	240	192902	5.00	ng		0.00
32) Perylene-d12	23.46	264	185515	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.25	112	1121	0.24	ng		0.00
5) Phenol-d6	6.80	99	1525	0.21	ng		0.00
7) Nitrobenzene-d5	8.76	82	1646	0.21	ng		0.00
13) 2,4,6-Tribromophenol	15.73	330	312	0.34	ng		0.00
14) 2-Fluorobiphenyl	12.86	172	3915	0.22	ng		0.00
27) Terphenyl-d14	19.59	244	6386	0.19	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.25	88	725	0.21	ng		97
3) n-Nitrosodimethylamine	3.54	42	927	0.20	ng	#	95
8) Nitrobenzene	8.80	77	1733	0.21	ng		98
9) Naphthalene	10.43	128	5051	0.22	ng		98
10) Hexachlorobutadiene	10.72	225	799	0.22	ng		98
11) 2-Methylnaphthalene	12.04	142	3261	0.20	ng		97
15) Acenaphthylene	13.95	152	22629	0.21	ng		100
16) Acenaphthene	14.30	154	3254	0.20	ng		91
17) Fluorene	15.29	166	4540	0.21	ng		98
19) 4-Bromophenyl-phenylether	16.18	248	1386	0.22	ng		95
20) Hexachlorobenzene	16.30	284	5737	0.21	ng		98
21) Pentachlorophenol	16.63	266	415	0.32	ng		91
22) Phenanthrene	17.01	178	8397	0.21	ng		99
23) Anthracene	17.10	178	7310	0.20	ng		99
24) Fluoranthene	19.01	202	8848	0.21	ng		100
26) Pyrene	19.38	202	9340	0.19	ng		100
28) Benzo(a)anthracene	21.12	228	9808	0.21	ng		99
29) Chrysene	21.16	228	10338	0.20	ng		98
30) Bis(2-ethylhexyl)phthalate	21.07	149	2078	0.36	ng		99
31) Indeno(1,2,3-cd)pyrene	25.86	276	9018	0.23	ng		100
33) Benzo(b)fluoranthene	22.75	252	7820	0.20	ng	#	98
34) Benzo(k)fluoranthene	22.79	252	9124	0.20	ng		96
35) Benzo(a)pyrene	23.35	252	7076	0.20	ng		96
36) Dibenzo(a,h)anthracene	25.89	278	7327	0.22	ng		96
37) Benzo(g,h,i)perylene	26.61	276	7408	0.21	ng		96

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

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