Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062315\
Data File : BE089929.D
Acq On : 24 Jun 2015 00:23
Operator : TP/IZ
Sample : SSTDICC0.2
Misc :
ALS Vial : 8 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.2

Quant Time: Jun 24 09:18:32 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 09:11:03 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	30473	5.00	ng		0.00
6) Naphthalene-d8	10.39	136	137961	5.00	ng		0.00
12) Acenaphthene-d10	14.25	164	86088	5.00	ng		0.00
18) Phenanthrene-d10	16.98	188	204391	5.00	ng		0.00
25) Chrysene-d12	21.14	240	227856	5.00	ng		0.00
32) Perylene-d12	23.47	264	203078	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.26	112	1057	0.15	ng		0.00
5) Phenol-d6	6.81	99	1604	0.18	ng		0.00
7) Nitrobenzene-d5	8.77	82	2070	0.20	ng		0.00
13) 2,4,6-Tribromophenol	15.74	330	167	0.07	ng		0.00
14) 2-Fluorobiphenyl	12.87	172	5278	0.18	ng		0.00
27) Terphenyl-d14	19.60	244	8075	0.24	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.26	88	973	0.28	ng		91
3) n-Nitrosodimethylamine	3.55	42	1131	0.21	ng	#	95
8) Nitrobenzene	8.81	77	2177	0.19	ng		98
9) Naphthalene	10.44	128	6580	0.20	ng		99
10) Hexachlorobutadiene	10.74	225	1076	0.19	ng		99
11) 2-Methylnaphthalene	12.05	142	4381	0.21	ng		97
15) Acenaphthylene	13.96	152	30946	0.58	ng		100
16) Acenaphthene	14.30	154	4618	0.18	ng		96
17) Fluorene	15.29	166	6400	0.29	ng	#	98
19) 4-Bromophenyl-phenylether	16.20	248	1769	0.21	ng		95
20) Hexachlorobenzene	16.30	284	8142	0.84	ng		99
21) Pentachlorophenol	16.65	266	215	0.09	ng	#	82
22) Phenanthrene	17.03	178	11276	0.17	ng		99
23) Anthracene	17.12	178	9679	0.31	ng		99
24) Fluoranthene	19.02	202	11476	0.25	ng		100
26) Pyrene	19.38	202	11806	0.15	ng		100
28) Benzo(a)anthracene	21.13	228	11171	0.17	ng		96
29) Chrysene	21.17	228	12668	0.18	ng		100
30) Bis(2-ethylhexyl)phthalate	21.07	149	971	0.03	ng	#	99
31) Indeno(1,2,3-cd)pyrene	25.89	276	7750	0.12	ng	#	100
33) Benzo(b)fluoranthene	22.76	252	8073	0.13	ng		97
34) Benzo(k)fluoranthene	22.81	252	9495	0.16	ng		99
35) Benzo(a)pyrene	23.37	252	6880	0.13	ng	#	95
36) Dibenzo(a,h)anthracene	25.91	278	6129	0.12	ng		97
37) Benzo(g,h,i)perylene	26.63	276	7061	0.13	ng		96

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

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