Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089986.D
Acq On : 29 Jun 2015 16:45
Operator : TP/IZ
Sample : SSTDICC0.8
Misc :
ALS Vial : 7 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC0.8

Quant Time: Jun 30 06:06:39 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	7.63	152	25148	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	106446	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	63863	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	156710	5.00	ng	0.00
25) Chrysene-d12	21.13	240	191926	5.00	ng	0.00
32) Perylene-d12	23.46	264	184965	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	4810	1.02	ng	0.00
5) Phenol-d6	6.80	99	6696	0.94	ng	0.00
7) Nitrobenzene-d5	8.76	82	7033	0.87	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	1510	1.56	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	16093	0.85	ng	0.00
27) Terphenyl-d14	19.59	244	27611	0.83	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	2901	0.84	ng	97
3) n-Nitrosodimethylamine	3.53	42	3925	0.86	ng	99
8) Nitrobenzene	8.80	77	7484	0.87	ng	100
9) Naphthalene	10.43	128	20227	0.85	ng	99
10) Hexachlorobutadiene	10.72	225	3208	0.85	ng	100
11) 2-Methylnaphthalene	12.05	142	13552	0.83	ng	96
15) Acenaphthylene	13.95	152	96538	0.86	ng	100
16) Acenaphthene	14.29	154	14074	0.85	ng	99
17) Fluorene	15.29	166	18849	0.84	ng	99
19) 4-Bromophenyl-phenylether	16.18	248	5629	0.87	ng	98
20) Hexachlorobenzene	16.30	284	23486	0.83	ng	99
21) Pentachlorophenol	16.63	266	2028	1.56	ng	97
22) Phenanthrene	17.01	178	33656	0.85	ng	100
23) Anthracene	17.10	178	31211	0.85	ng	100
24) Fluoranthene	19.02	202	37520	0.87	ng	99
26) Pyrene	19.38	202	39247	0.81	ng	100
28) Benzo(a)anthracene	21.12	228	40452	0.87	ng	99
29) Chrysene	21.16	228	42099	0.83	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	11275	1.98	ng	100
31) Indeno(1,2,3-cd)pyrene	25.86	276	40426	1.05	ng	100
33) Benzo(b)fluoranthene	22.75	252	34811	0.88	ng	99
34) Benzo(k)fluoranthene	22.79	252	38735	0.86	ng	100
35) Benzo(a)pyrene	23.35	252	31995	0.90	ng	99
36) Dibenzo(a,h)anthracene	25.89	278	32987	0.99	ng	100
37) Benzo(g,h,i)perylene	26.61	276	33395	0.94	ng	99

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

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