Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089984.D
Acq On : 29 Jun 2015 15:33
Operator : TP/IZ
Sample : SSTDICC002
Misc :
ALS Vial : 5 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDICC002

Quant Time: Jun 30 06:05:59 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	27942	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	121318	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	72076	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	177913	5.00	ng	0.00
25) Chrysene-d12	21.13	240	229780	5.00	ng	0.00
32) Perylene-d12	23.46	264	223505	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	11728	2.23	ng	0.00
5) Phenol-d6	6.80	99	16748	2.11	ng	0.00
7) Nitrobenzene-d5	8.76	82	17449	1.90	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	4165	3.82	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	38165	1.79	ng	0.00
27) Terphenyl-d14	19.59	244	69446	1.74	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	6556	1.71	ng	99
3) n-Nitrosodimethylamine	3.53	42	9161	1.81	ng	98
8) Nitrobenzene	8.80	77	18741	1.92	ng	99
9) Naphthalene	10.43	128	47800	1.75	ng	100
10) Hexachlorobutadiene	10.72	225	7423	1.73	ng	99
11) 2-Methylnaphthalene	12.04	142	32258	1.73	ng	99
15) Acenaphthylene	13.95	152	233589	1.83	ng	100
16) Acenaphthene	14.30	154	33979	1.82	ng	98
17) Fluorene	15.29	166	44259	1.74	ng	99
19) 4-Bromophenyl-phenylether	16.18	248	13395	1.81	ng	100
20) Hexachlorobenzene	16.30	284	54277	1.70	ng	99
21) Pentachlorophenol	16.63	266	6186	4.20	ng	97
22) Phenanthrene	17.01	178	79943	1.77	ng	100
23) Anthracene	17.10	178	76015	1.82	ng	99
24) Fluoranthene	19.02	202	94558	1.92	ng	100
26) Pyrene	19.38	202	98795	1.70	ng	100
28) Benzo(a)anthracene	21.11	228	102360	1.84	ng	99
29) Chrysene	21.17	228	106600	1.76	ng	99
30) Bis(2-ethylhexyl)phthalate	21.07	149	39142	5.74	ng	100
31) Indeno(1,2,3-cd)pyrene	25.87	276	107540	2.34	ng	99
33) Benzo(b)fluoranthene	22.75	252	91224	1.91	ng	99
34) Benzo(k)fluoranthene	22.80	252	101784	1.88	ng	99
35) Benzo(a)pyrene	23.35	252	86310	2.02	ng	99
36) Dibenzo(a,h)anthracene	25.89	278	87783	2.18	ng	99
37) Benzo(g,h,i)perylene	26.60	276	88004	2.05	ng	99

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

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