Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
Data File : BE089990.D
Acq On : 29 Jun 2015 19:08
Operator : TP/IZ
Sample : SSTDICV001
Misc :
ALS Vial : 11 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : ICVBE062915

Quant Time: Jun 30 07:02:35 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 07:01:01 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	26423	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	115038	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	70141	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	173058	5.00	ng	0.00
25) Chrysene-d12	21.13	240	218956	5.00	ng	0.00
32) Perylene-d12	23.45	264	212785	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	5077	0.84	ng	0.00
5) Phenol-d6	6.80	99	7318	0.86	ng	0.00
7) Nitrobenzene-d5	8.76	82	7678	0.84	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	1595	0.78	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	18616	0.88	ng	0.00
27) Terphenyl-d14	19.59	244	32261	0.88	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.24	88	3051	0.84	ng	99
3) n-Nitrosodimethylamine	3.53	42	4123	0.84	ng	99
8) Nitrobenzene	8.80	77	8162	0.84	ng	99
9) Naphthalene	10.43	128	23683	0.91	ng	99
10) Hexachlorobutadiene	10.72	225	3787	0.93	ng	99
11) 2-Methylnaphthalene	12.04	142	15579	0.90	ng	100
15) Acenaphthylene	13.95	152	114139	0.90	ng	100
16) Acenaphthene	14.30	154	17169	0.94	ng	97
17) Fluorene	15.29	166	22865	0.92	ng	100
19) 4-Bromophenyl-phenylether	16.18	248	6542	0.89	ng	98
20) Hexachlorobenzene	16.30	284	28416	0.93	ng	98
21) Pentachlorophenol	16.63	266	2008	0.75	ng	97
22) Phenanthrene	17.01	178	40858	0.92	ng	100
23) Anthracene	17.10	178	37448	0.92	ng	99
24) Fluoranthene	19.01	202	45676	0.92	ng	100
26) Pyrene	19.37	202	47736	0.91	ng	100
28) Benzo(a)anthracene	21.11	228	50139	0.92	ng	99
29) Chrysene	21.16	228	50803	0.89	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	10708	0.71	ng	99
31) Indeno(1,2,3-cd)pyrene	25.87	276	49677	0.91	ng	100
33) Benzo(b)fluoranthene	22.74	252	44517	0.95	ng	99
34) Benzo(k)fluoranthene	22.79	252	47844	0.91	ng	99
35) Benzo(a)pyrene	23.35	252	39358	0.91	ng	99
36) Dibenzo(a,h)anthracene	25.89	278	39902	0.90	ng	99
37) Benzo(g,h,i)perylene	26.60	276	40513	0.90	ng	100

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

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