Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090115\
Data File : BE090688.D
Acq On : 2 Sep 2015 4:25
Operator : UM/IZ
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDCCC001

Manual Integrations APPROVED
MMDadoda
9/2/2015 4:08:21 PM

Quant Time: Sep 02 12:25:22 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Wed Sep 02 11:40:50 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	43351	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	188167	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	90978	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	221655	5.00	ng	0.00
25) Chrysene-d12	21.07	240	300813	5.00	ng	0.00
32) Perylene-d12	23.36	264	264131	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	8279	1.00	ng	0.00
5) Phenol-d6	6.76	99	11428	0.95	ng	0.00
7) Nitrobenzene-d5	8.71	82	11991	0.96	ng	0.00
13) 2,4,6-Tribromophenol	15.67	330	1980	0.88	ng	0.00
14) 2-Fluorobiphenyl	12.80	172	24862	0.99	ng	0.00
27) Terphenyl-d14	19.54	244	31278	0.94	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.16	88	7083	1.12	ng	96
3) n-Nitrosodimethylamine	3.46	42	8168	1.00	ng	96
8) Nitrobenzene	8.75	77	12763	0.93	ng	98
9) Naphthalene	10.35	128	41100	1.00	ng	99
10) Hexachlorobutadiene	10.65	225	6534	1.02	ng	100
11) 2-Methylnaphthalene	11.99	142	21840	0.98	ng	99
15) Acenaphthylene	13.89	152	142800	0.98	ng	100
16) Acenaphthene	14.23	154	24789	1.02	ng	99
17) Fluorene	15.22	166	29613	0.98	ng	98
19) 4-Bromophenyl-phenylether	16.13	248	7456	0.99	ng	98
20) Hexachlorobenzene	16.23	284	41830	1.06	ng	97
21) Pentachlorophenol	16.60	266	1816m	1.02	ng	
22) Phenanthrene	16.96	178	55561	1.01	ng	99
23) Anthracene	17.05	178	44782	0.98	ng	99
24) Fluoranthene	18.96	202	55978	0.97	ng	100
26) Pyrene	19.32	202	59514	0.96	ng	100
28) Benzo(a)anthracene	21.06	228	67439	1.01	ng	99
29) Chrysene	21.11	228	76878	1.05	ng	100
30) Bis(2-ethylhexyl)phthalate	21.01	149	13298	0.80	ng	99
31) Indeno(1,2,3-cd)pyrene	25.72	276	65798	0.93	ng	100
33) Benzo(b)fluoranthene	22.66	252	55655	0.97	ng	98
34) Benzo(k)fluoranthene	22.71	252	68034	1.02	ng	100
35) Benzo(a)pyrene	23.26	252	49359	0.96	ng	99
36) Dibenzo(a,h)anthracene	25.74	278	51535	0.94	ng	98
37) Benzo(g,h,i)perylene	26.45	276	55563	0.94	ng	99

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

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