Quantitation Report (QT Reviewed)

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090342.D
Acq On : 4 Aug 2015 11:50
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument : BNA_E
ClientSampleId : SSTDCCC001EC

Quant Time: Aug 04 13:22:23 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 2015
Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units		Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2951	5.00	ng		0.00
6) Naphthalene-d8	10.35	136	13097	5.00	ng		0.00
12) Acenaphthene-d10	14.21	164	5515	5.00	ng		0.00
18) Phenanthrene-d10	16.94	188	10213	5.00	ng		0.00
25) Chrysene-d12	21.11	240	7588	5.00	ng		0.00
32) Perylene-d12	23.42	264	6763	5.00	ng		0.00
System Monitoring Compounds							
4) 2-Fluorophenol	5.22	112	796	1.41	ng		0.00
5) Phenol-d6	6.78	99	906	1.31	ng		0.00
7) Nitrobenzene-d5	8.74	82	1130	1.46	ng		0.00
13) 2,4,6-Tribromophenol	15.73	330	146	2.00	ng		0.00
14) 2-Fluorobiphenyl	12.84	172	2828	1.77	ng		0.00
27) Terphenyl-d14	19.58	244	2722	1.92	ng		0.00
Target Compounds							Qvalue
2) 1,4-Dioxane	3.20	88	733	2.01	ng	#	72
3) n-Nitrosodimethylamine	3.51	42	520	1.04	ng	#	95
8) Nitrobenzene	8.78	77	643	0.81	ng		98
9) Naphthalene	10.40	128	2319	0.82	ng		94
10) Hexachlorobutadiene	10.69	225	317	0.84	ng		98
11) 2-Methylnaphthalene	12.02	142	1316	0.76	ng		97
15) Acenaphthylene	13.93	152	7899	0.82	ng		100
16) Acenaphthene	14.27	154	1071	0.77	ng		96
17) Fluorene	15.26	166	1238	0.73	ng	#	98
19) 4-Bromophenyl-phenylether	16.16	248	335	0.87	ng		95
20) Hexachlorobenzene	16.27	284	1265	0.87	ng		99
21) Pentachlorophenol	16.63	266	69	1.10	ng		95
22) Phenanthrene	17.00	178	1807	0.78	ng		99
23) Anthracene	17.08	178	1544	0.74	ng		96
24) Fluoranthene	19.00	202	1774	0.76	ng		100
26) Pyrene	19.36	202	1776	0.87	ng		99
28) Benzo(a)anthracene	21.10	228	1004	0.70	ng		94
29) Chrysene	21.15	228	1586	0.87	ng		94
30) Bis(2-ethylhexyl)phthalate	21.04	149	429	0.99	ng		95
31) Indeno(1,2,3-cd)pyrene	25.83	276	913	1.18	ng	#	91
33) Benzo(b)fluoranthene	22.72	252	835	0.95	ng	#	90
34) Benzo(k)fluoranthene	22.76	252	1316	0.72	ng	#	86
35) Benzo(a)pyrene	23.32	252	822	0.95	ng	#	83
36) Dibenzo(a,h)anthracene	25.85	278	475	0.93	ng	#	88
37) Benzo(g,h,i)perylene	26.55	276	995	1.00	ng	#	91

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

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
Data File : BE090342.D
Acq On : 4 Aug 2015 11:50
Operator : TP/UM
Sample : SSTDCCC001
Misc :
ALS Vial : 6 Sample Multiplier: 1

Instrument :
BNA_E
ClientSampleId :
SSTDCCC001EC

Quant Time: Aug 04 13:22:23 2015
Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Tue Aug 04 02:31:31 2015
Response via : Initial Calibration