

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071015\
 Data File : BE090081.D
 Acq On : 10 Jul 2015 12:37
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Jul 10 23:08:10 2015
 Quant Method : Z:\HPCHEM1\BNA_E\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	7.61	152	24410	5.00	ng	-0.02
6) Naphthalene-d8	10.36	136	103486	5.00	ng	-0.02
12) Acenaphthene-d10	14.22	164	56356	5.00	ng	-0.01
18) Phenanthrene-d10	16.96	188	123624	5.00	ng	-0.02
25) Chrysene-d12	21.12	240	134419	5.00	ng	-0.01
32) Perylene-d12	23.44	264	123194	5.00	ng	-0.02
System Monitoring Compounds						
4) 2-Fluorophenol	5.24	112	5159	0.92	ng	-0.01
5) Phenol-d6	6.79	99	7439	0.95	ng	0.00
7) Nitrobenzene-d5	8.74	82	7198	0.88	ng	-0.02
13) 2,4,6-Tribromophenol	15.73	330	1382	0.84	ng	0.00
14) 2-Fluorobiphenyl	12.84	172	15567	0.92	ng	-0.02
27) Terphenyl-d14	19.58	244	21013	0.94	ng	-0.01
Target Compounds						Qvalue
2) 1,4-Dioxane	3.23	88	2991	0.89	ng	90
3) n-Nitrosodimethylamine	3.51	42	3700	0.81	ng	# 93
8) Nitrobenzene	8.78	77	7581	0.87	ng	99
9) Naphthalene	10.41	128	21587	0.92	ng	99
10) Hexachlorobutadiene	10.71	225	3514	0.96	ng	100
11) 2-Methylnaphthalene	12.03	142	13587	0.87	ng	99
15) Acenaphthylene	13.93	152	93336	0.92	ng	100
16) Acenaphthene	14.28	154	13610	0.93	ng	98
17) Fluorene	15.27	166	17828	0.89	ng	98
19) 4-Bromophenyl-phenylether	16.16	248	4774	0.91	ng	97
20) Hexachlorobenzene	16.28	284	19698	0.90	ng	99
21) Pentachlorophenol	16.63	266	11006	4.06	ng	98
22) Phenanthrene	17.00	178	27633	0.87	ng	100
23) Anthracene	17.10	178	25851	0.89	ng	99
24) Fluoranthene	19.00	202	30303	0.86	ng	100
26) Pyrene	19.36	202	31771	0.99	ng	100
28) Benzo(a)anthracene	21.10	228	30990	0.92	ng	100
29) Chrysene	21.16	228	31440	0.90	ng	99
30) Bis(2-ethylhexyl)phthalate	21.05	149	11272	0.99	ng	98
31) Indeno(1,2,3-cd)pyrene	25.84	276	31498	0.94	ng	99
33) Benzo(b)fluoranthene	22.73	252	27412	1.01	ng	99
34) Benzo(k)fluoranthene	22.77	252	30787	1.01	ng	99
35) Benzo(a)pyrene	23.33	252	25505	1.02	ng	99
36) Dibenzo(a,h)anthracene	25.85	278	24969	0.97	ng	98
37) Benzo(g,h,i)perylene	26.57	276	25562	0.98	ng	98

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

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