

Data Path : Z:\HPCHEM1\BNA E\DATA\BE081415\
 Data File : BE090514.D
 Acq On : 14 Aug 2015 10:06
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Aug 14 15:26:29 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\8270-SIM-BE081415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 14 15:17:14 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	24995	5.00	ng	0.00
6) Naphthalene-d8	10.33	136	124435	5.00	ng	0.00
12) Acenaphthene-d10	14.19	164	69430	5.00	ng	0.00
18) Phenanthrene-d10	16.94	188	155622	5.00	ng	0.00
25) Chrysene-d12	21.09	240	160497	5.00	ng	0.00
32) Perylene-d12	23.40	264	147016	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.21	112	7341	0.96	ng	0.00
5) Phenol-d6	6.77	99	10760	0.99	ng	0.00
7) Nitrobenzene-d5	8.72	82	9944	1.01	ng	0.00
13) 2,4,6-Tribromophenol	15.69	330	2657	1.02	ng	-0.02
14) 2-Fluorobiphenyl	12.81	172	20838	0.96	ng	0.00
27) Terphenyl-d14	19.55	244	29268	0.95	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.19	88	3426	0.95	ng	97
3) n-Nitrosodimethylamine	3.48	42	4541	1.01	ng	# 77
8) Nitrobenzene	8.76	77	10308	1.02	ng	97
9) Naphthalene	10.38	128	28230	0.94	ng	99
10) Hexachlorobutadiene	10.68	225	4128	0.93	ng	99
11) 2-Methylnaphthalene	12.00	142	18895	0.98	ng	96
15) Acenaphthylene	13.91	152	127444	0.97	ng	100
16) Acenaphthene	14.25	154	18965	0.94	ng	# 79
17) Fluorene	15.24	166	23661	0.96	ng	99
19) 4-Bromophenyl-phenylether	16.14	248	6476	0.96	ng	86
20) Hexachlorobenzene	16.27	284	27689	0.92	ng	94
21) Pentachlorophenol	16.61	266	2317	1.11	ng	91
22) Phenanthrene	16.98	178	39909	0.92	ng	100
23) Anthracene	17.07	178	37769	0.97	ng	100
24) Fluoranthene	18.98	202	44363	0.98	ng	99
26) Pyrene	19.34	202	45240	0.94	ng	99
28) Benzo(a)anthracene	21.08	228	41104	0.92	ng	98
29) Chrysene	21.13	228	43391	0.95	ng	98
30) Bis(2-ethylhexyl)phthalate	21.03	149	24063	1.06	ng	99
31) Indeno(1,2,3-cd)pyrene	25.78	276	39950	0.97	ng	# 92
33) Benzo(b)fluoranthene	22.70	252	35582	0.98	ng	99
34) Benzo(k)fluoranthene	22.74	252	39843	0.96	ng	98
35) Benzo(a)pyrene	23.29	252	32742	0.99	ng	# 97
36) Dibenzo(a,h)anthracene	25.80	278	31092	0.98	ng	96
37) Benzo(g,h,i)perylene	26.51	276	34362	0.98	ng	96

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

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