

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
 Data File : BE089997.D
 Acq On : 30 Jun 2015 1:09
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Jun 30 07:17:49 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	7.63	152	18038	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	75914	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	45351	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	111411	5.00	ng	0.00
25) Chrysene-d12	21.13	240	140177	5.00	ng	0.00
32) Perylene-d12	23.46	264	129405	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	3253	0.79	ng	0.00
5) Phenol-d6	6.80	99	4524	0.78	ng	0.00
7) Nitrobenzene-d5	8.76	82	5226	0.87	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	799	0.60	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	13258	0.97	ng	0.00
27) Terphenyl-d14	19.59	244	22260	0.95	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.24	88	2415	0.98	ng	96
3) n-Nitrosodimethylamine	3.53	42	3047	0.91	ng	# 92
8) Nitrobenzene	8.80	77	5589	0.87	ng	98
9) Naphthalene	10.43	128	16485	0.96	ng	100
10) Hexachlorobutadiene	10.72	225	2753	1.02	ng	99
11) 2-Methylnaphthalene	12.04	142	10883	0.95	ng	100
15) Acenaphthylene	13.95	152	77415	0.95	ng	100
16) Acenaphthene	14.29	154	11459	0.97	ng	96
17) Fluorene	15.29	166	15376	0.96	ng	99
19) 4-Bromophenyl-phenylether	16.18	248	4580	0.97	ng	100
20) Hexachlorobenzene	16.30	284	19763	1.00	ng	99
21) Pentachlorophenol	16.63	266	1056	0.66	ng	97
22) Phenanthrene	17.01	178	27473	0.96	ng	100
23) Anthracene	17.10	178	25138	0.96	ng	100
24) Fluoranthene	19.01	202	30362	0.95	ng	100
26) Pyrene	19.37	202	31747	0.95	ng	100
28) Benzo(a)anthracene	21.11	228	32587	0.93	ng	99
29) Chrysene	21.16	228	35065	0.96	ng	98
30) Bis(2-ethylhexyl)phthalate	21.06	149	4826	0.59	ng	100
31) Indeno(1,2,3-cd)pyrene	25.86	276	29428	0.84	ng	100
33) Benzo(b)fluoranthene	22.74	252	27580	0.97	ng	99
34) Benzo(k)fluoranthene	22.79	252	30819	0.97	ng	99
35) Benzo(a)pyrene	23.35	252	24069	0.91	ng	100
36) Dibenzo(a,h)anthracene	25.89	278	23992	0.89	ng	100
37) Benzo(g,h,i)perylene	26.60	276	24706	0.90	ng	99

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

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