

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062315\
 Data File : BE089926.D
 Acq On : 23 Jun 2015 22:35
 Operator : TP/IZ
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICCC001

Quant Time: Jun 24 09:17:12 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 24 09:11:03 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	29381	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	134558	5.00	ng	0.00
12) Acenaphthene-d10	14.25	164	84165	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	207721	5.00	ng	0.00
25) Chrysene-d12	21.14	240	222554	5.00	ng	0.00
32) Perylene-d12	23.47	264	195570	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	4991	0.75	ng	0.00
5) Phenol-d6	6.81	99	7531	0.87	ng	0.00
7) Nitrobenzene-d5	8.77	82	8961	0.87	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	902	0.38	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	22032	0.76	ng	0.00
27) Terphenyl-d14	19.60	244	34925	1.05	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.26	88	3377	1.01	ng	99
3) n-Nitrosodimethylamine	3.55	42	4828	0.94	ng	99
8) Nitrobenzene	8.81	77	9481	0.86	ng	100
9) Naphthalene	10.45	128	27706	0.86	ng	100
10) Hexachlorobutadiene	10.74	225	4396	0.81	ng	100
11) 2-Methylnaphthalene	12.05	142	18622	0.90	ng	100
15) Acenaphthylene	13.96	152	137498	2.63	ng	100
16) Acenaphthene	14.30	154	20506	0.84	ng	100
17) Fluorene	15.29	166	27249	1.24	ng	100
19) 4-Bromophenyl-phenylether	16.20	248	7393	0.87	ng	99
20) Hexachlorobenzene	16.30	284	33402	3.39	ng	100
21) Pentachlorophenol	16.65	266	1062	0.44	ng	100
22) Phenanthrene	17.03	178	46745	0.68	ng	100
23) Anthracene	17.12	178	43761	1.38	ng	100
24) Fluoranthene	19.02	202	51343	1.11	ng	100
26) Pyrene	19.38	202	53506	0.71	ng	100
28) Benzo(a)anthracene	21.12	228	48684	0.75	ng	100
29) Chrysene	21.17	228	52240	0.76	ng	100
30) Bis(2-ethylhexyl)phthalate	21.07	149	3820	0.11	ng	100
31) Indeno(1,2,3-cd)pyrene	25.89	276	40406	0.63	ng	# 100
33) Benzo(b)fluoranthene	22.76	252	39169	0.64	ng	100
34) Benzo(k)fluoranthene	22.81	252	44057	0.78	ng	100
35) Benzo(a)pyrene	23.36	252	34130	0.66	ng	100
36) Dibenzo(a,h)anthracene	25.91	278	31704	0.62	ng	100
37) Benzo(g,h,i)perylene	26.62	276	34438	0.64	ng	100

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

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