

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089979.D
 Acq On : 27 Jun 2015 16:31
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Jun 29 00:57:28 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\8270-SIM-BE062315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 25 16:51:27 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	29511	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	145581	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	86631	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	193145	5.00	ng	0.00
25) Chrysene-d12	21.14	240	206236	5.00	ng	0.00
32) Perylene-d12	23.47	264	191778	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	8657	1.62	ng	0.00
5) Phenol-d6	6.81	99	17006	2.08	ng	0.00
7) Nitrobenzene-d5	8.77	82	13063	1.21	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	3268	2.37	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	32131	1.27	ng	0.00
27) Terphenyl-d14	19.60	244	45809	1.29	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.26	88	3360	0.94	ng	95
3) n-Nitrosodimethylamine	3.55	42	4556	0.86	ng	# 92
8) Nitrobenzene	8.81	77	9667	0.84	ng	99
9) Naphthalene	10.44	128	27151	0.84	ng	100
10) Hexachlorobutadiene	10.73	225	4193	0.82	ng	100
11) 2-Methylnaphthalene	12.05	142	17704	0.80	ng	99
15) Acenaphthylene	13.95	152	125490	0.83	ng	100
16) Acenaphthene	14.30	154	18086	0.81	ng	98
17) Fluorene	15.29	166	23216	0.83	ng	98
19) 4-Bromophenyl-phenylether	16.18	248	6225	0.79	ng	96
20) Hexachlorobenzene	16.30	284	28057	0.81	ng	99
21) Pentachlorophenol	16.63	266	1967	1.42	ng	# 88
22) Phenanthrene	17.01	178	39797	0.82	ng	99
23) Anthracene	17.12	178	35768	0.80	ng	99
24) Fluoranthene	19.02	202	44747	0.85	ng	100
26) Pyrene	19.39	202	46364	0.89	ng	100
28) Benzo(a)anthracene	21.12	228	43685	0.89	ng	99
29) Chrysene	21.17	228	42081	0.78	ng	98
30) Bis(2-ethylhexyl)phthalate	21.07	149	11230	2.67	ng	99
31) Indeno(1,2,3-cd)pyrene	25.88	276	43632	1.10	ng	# 100
33) Benzo(b)fluoranthene	22.76	252	39459	0.97	ng	99
34) Benzo(k)fluoranthene	22.80	252	40042	0.87	ng	99
35) Benzo(a)pyrene	23.36	252	35614	0.99	ng	100
36) Dibenzo(a,h)anthracene	25.90	278	33133	0.99	ng	99
37) Benzo(g,h,i)perylene	26.62	276	37025	1.03	ng	99

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

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