

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE090415\
 Data File : BE090721.D
 Acq On : 4 Sep 2015 9:46
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Sep 04 13:13:17 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE090115.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Sep 02 11:40:50 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	38149	5.00	ng	0.00
6) Naphthalene-d8	10.31	136	176800	5.00	ng	0.00
12) Acenaphthene-d10	14.17	164	97866	5.00	ng	0.00
18) Phenanthrene-d10	16.91	188	246703	5.00	ng	0.00
25) Chrysene-d12	21.07	240	325536	5.00	ng	0.00
32) Perylene-d12	23.36	264	308715	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.19	112	7425m	1.01	ng	0.00
5) Phenol-d6	6.76	99	11451	1.08	ng	0.01
7) Nitrobenzene-d5	8.71	82	11411	0.98	ng	0.00
13) 2,4,6-Tribromophenol	15.67	330	3103	1.22	ng	0.00
14) 2-Fluorobiphenyl	12.80	172	26412	0.97	ng	0.00
27) Terphenyl-d14	19.54	244	39288	1.09	ng	0.00

Target Compounds					Qvalue	
2) 1,4-Dioxane	3.15	88	5841	1.03	ng	98
3) n-Nitrosodimethylamine	3.45	42	6632	0.91	ng	85
8) Nitrobenzene	8.75	77	12200	0.94	ng	99
9) Naphthalene	10.35	128	38538	1.00	ng	99
10) Hexachlorobutadiene	10.65	225	5792	0.96	ng	99
11) 2-Methylnaphthalene	11.99	142	23281	1.12	ng	99
15) Acenaphthylene	13.89	152	163761	1.05	ng	100
16) Acenaphthene	14.23	154	26610	1.02	ng	93
17) Fluorene	15.22	166	33902	1.04	ng	97
19) 4-Bromophenyl-phenylether	16.13	248	9038	1.07	ng	92
20) Hexachlorobenzene	16.23	284	45513	1.03	ng	96
21) Pentachlorophenol	16.60	266	2518	1.15	ng	94
22) Phenanthrene	16.94	178	63902	1.05	ng	99
23) Anthracene	17.05	178	56466	1.11	ng	100
24) Fluoranthene	18.96	202	72054	1.12	ng	99
26) Pyrene	19.32	202	76386	1.13	ng	100
28) Benzo(a)anthracene	21.05	228	77320	1.07	ng	99
29) Chrysene	21.11	228	88163	1.12	ng	99
30) Bis(2-ethylhexyl)phthalate	21.00	149	24636	1.28	ng	100
31) Indeno(1,2,3-cd)pyrene	25.71	276	88209	1.15	ng	99
33) Benzo(b)fluoranthene	22.66	252	70803	1.06	ng	100
34) Benzo(k)fluoranthene	22.71	252	81937	1.05	ng	99
35) Benzo(a)pyrene	23.26	252	66287	1.11	ng	99
36) Dibenzo(a,h)anthracene	25.74	278	68565	1.05	ng	99
37) Benzo(g,h,i)perylene	26.44	276	72811	1.06	ng	97

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

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