

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062915\
 Data File : BE089981.D
 Acq On : 29 Jun 2015 13:19
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001

Quant Time: Jun 30 06:04:53 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 06:02:58 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	31499	5.00	ng	0.00
6) Naphthalene-d8	10.38	136	138199	5.00	ng	0.00
12) Acenaphthene-d10	14.23	164	86283	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	210841	5.00	ng	0.00
25) Chrysene-d12	21.13	240	267006	5.00	ng	0.00
32) Perylene-d12	23.46	264	265950	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.25	112	7328	1.24	ng	0.00
5) Phenol-d6	6.80	99	10409	1.16	ng	0.00
7) Nitrobenzene-d5	8.76	82	9961	0.95	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	2832	2.17	ng	0.00
14) 2-Fluorobiphenyl	12.86	172	22811	0.90	ng	0.00
27) Terphenyl-d14	19.59	244	42153	0.91	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.25	88	3899	0.90	ng	95
3) n-Nitrosodimethylamine	3.53	42	5267	0.92	ng	# 97
8) Nitrobenzene	8.80	77	10423	0.94	ng	99
9) Naphthalene	10.43	128	28714	0.93	ng	99
10) Hexachlorobutadiene	10.72	225	4384	0.89	ng	99
11) 2-Methylnaphthalene	12.05	142	19424	0.92	ng	96
15) Acenaphthylene	13.95	152	144775	0.95	ng	100
16) Acenaphthene	14.30	154	21493	0.96	ng	98
17) Fluorene	15.29	166	28238	0.93	ng	99
19) 4-Bromophenyl-phenylether	16.18	248	8252	0.94	ng	98
20) Hexachlorobenzene	16.30	284	34064	0.90	ng	99
21) Pentachlorophenol	16.63	266	4236	2.43	ng	97
22) Phenanthrene	17.01	178	50921	0.95	ng	99
23) Anthracene	17.10	178	48116	0.97	ng	100
24) Fluoranthene	19.02	202	59605	1.02	ng	100
26) Pyrene	19.38	202	62266	0.92	ng	100
28) Benzo(a)anthracene	21.11	228	64368	1.00	ng	100
29) Chrysene	21.16	228	63061	0.89	ng	97
30) Bis(2-ethylhexyl)phthalate	21.07	149	26623	3.36	ng	100
31) Indeno(1,2,3-cd)pyrene	25.87	276	69136	1.30	ng	100
33) Benzo(b)fluoranthene	22.75	252	58653	1.03	ng	100
34) Benzo(k)fluoranthene	22.79	252	62728	0.97	ng	100
35) Benzo(a)pyrene	23.35	252	54050	1.06	ng	99
36) Dibenzo(a,h)anthracene	25.89	278	55561	1.16	ng	100
37) Benzo(g,h,i)perylene	26.60	276	56049	1.10	ng	99

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

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