

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE062615\
 Data File : BE089953.D
 Acq On : 26 Jun 2015 21:25
 Operator : TP/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDCCC001EC

Quant Time: Jun 27 01:01:53 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	25354	5.00	ng	0.00
6) Naphthalene-d8	10.39	136	118414	5.00	ng	0.00
12) Acenaphthene-d10	14.24	164	73983	5.00	ng	0.00
18) Phenanthrene-d10	16.98	188	181271	5.00	ng	0.00
25) Chrysene-d12	21.13	240	210592	5.00	ng	0.00
32) Perylene-d12	23.46	264	185978	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	5.26	112	4900	1.07	ng	0.00
5) Phenol-d6	6.81	99	11250	1.60	ng	0.00
7) Nitrobenzene-d5	8.77	82	9394	1.07	ng	0.00
13) 2,4,6-Tribromophenol	15.74	330	1701	1.57	ng	0.00
14) 2-Fluorobiphenyl	12.87	172	27974	1.30	ng	0.00
27) Terphenyl-d14	19.60	244	45885	1.26	ng	0.00

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.26	88	3130	1.02	ng	96
3) n-Nitrosodimethylamine	3.54	42	4136	0.91	ng	# 96
8) Nitrobenzene	8.81	77	9853	1.05	ng	99
9) Naphthalene	10.44	128	28900	1.10	ng	100
10) Hexachlorobutadiene	10.74	225	4144	1.00	ng	99
11) 2-Methylnaphthalene	12.05	142	22330	1.24	ng	98
15) Acenaphthylene	13.96	152	166648	1.28	ng	100
16) Acenaphthene	14.30	154	25024	1.31	ng	99
17) Fluorene	15.29	166	33686	1.41	ng	98
19) 4-Bromophenyl-phenylether	16.18	248	9269	1.25	ng	98
20) Hexachlorobenzene	16.30	284	41251	1.27	ng	98
21) Pentachlorophenol	16.63	266	2242	1.63	ng	91
22) Phenanthrene	17.01	178	57276	1.26	ng	99
23) Anthracene	17.12	178	54293	1.29	ng	99
24) Fluoranthene	19.02	202	64528	1.31	ng	100
26) Pyrene	19.38	202	67512	1.27	ng	99
28) Benzo(a)anthracene	21.12	228	67111	1.33	ng	98
29) Chrysene	21.17	228	70595	1.28	ng	99
30) Bis(2-ethylhexyl)phthalate	21.07	149	6585	1.53	ng	100
31) Indeno(1,2,3-cd)pyrene	25.88	276	58095	1.44	ng	# 100
33) Benzo(b)fluoranthene	22.75	252	54617	1.39	ng	99
34) Benzo(k)fluoranthene	22.80	252	62359	1.40	ng	99
35) Benzo(a)pyrene	23.36	252	49159	1.41	ng	100
36) Dibenzo(a,h)anthracene	25.90	278	46871	1.44	ng	99
37) Benzo(g,h,i)perylene	26.62	276	49432	1.41	ng	99

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

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