

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE080315\
 Data File : BE090321.D
 Acq On : 3 Aug 2015 22:25
 Operator : TP/UM
 Sample : SSTDICCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTDICCC001

Manual Integrations
 APPROVED

MMDadoda
 8/4/2015 3:26:49 PM

Quant Time: Aug 04 02:15:08 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\8270-SIM-BE080315.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 04 01:38:34 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	6236	5.00	ng	0.00
6) Naphthalene-d8	10.36	136	24930	5.00	ng	0.00
12) Acenaphthene-d10	14.21	164	12621	5.00	ng	0.00
18) Phenanthrene-d10	16.95	188	26451	5.00	ng	0.00
25) Chrysene-d12	21.11	240	20918	5.00	ng	0.00
32) Perylene-d12	23.43	264	16784	5.00	ng	0.00
System Monitoring Compounds						
4) 2-Fluorophenol	5.22	112	1158	0.96	ng	0.00
5) Phenol-d6	6.78	99	1359	0.96	ng	0.00
7) Nitrobenzene-d5	8.74	82	1418	0.93	ng	0.00
13) 2,4,6-Tribromophenol	15.73	330	121	0.81	ng	0.00
14) 2-Fluorobiphenyl	12.83	172	3696	1.02	ng	0.00
27) Terphenyl-d14	19.58	244	4095	1.03	ng	0.00
Target Compounds						Qvalue
2) 1,4-Dioxane	3.21	88	1003	1.09	ng	100
3) n-Nitrosodimethylamine	3.51	42	1068	1.00	ng	100
8) Nitrobenzene	8.78	77	1433	0.94	ng	100
9) Naphthalene	10.40	128	5317	1.00	ng	100
10) Hexachlorobutadiene	10.69	225	735	1.01	ng	100
11) 2-Methylnaphthalene	12.02	142	3283	0.99	ng	100
15) Acenaphthylene	13.93	152	21917	1.00	ng	100
16) Acenaphthene	14.27	154	3161	1.00	ng	100
17) Fluorene	15.26	166	3623	0.95	ng	100
19) 4-Bromophenyl-phenylether	16.16	248	1022	1.02	ng	100
20) Hexachlorobenzene	16.27	284	3839	1.02	ng	100
21) Pentachlorophenol	16.63	266	130	0.75	ng	100
22) Phenanthrene	17.00	178	5776	0.96	ng	100
23) Anthracene	17.08	178	5076	0.95	ng	100
24) Fluoranthene	19.00	202	5810	0.97	ng	100
26) Pyrene	19.36	202	5937	0.97	ng	100
28) Benzo(a)anthracene	21.10	228	3554	1.07	ng	100
29) Chrysene	21.15	228	5777m	0.93	ng	
30) Bis(2-ethylhexyl)phthalate	21.04	149	994	0.78	ng	100
31) Indeno(1,2,3-cd)pyrene	25.83	276	1645	1.22	ng	100
33) Benzo(b)fluoranthene	22.72	252	1948	1.28	ng	100
34) Benzo(k)fluoranthene	22.76	252	4880m	1.21	ng	
35) Benzo(a)pyrene	23.32	252	1996	1.14	ng	100
36) Dibenzo(a,h)anthracene	25.86	278	1017	1.17	ng	100
37) Benzo(g,h,i)perylene	26.56	276	2421	1.30	ng	100

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

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