

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM052215\
 Data File : BM001392.D
 Acq On : 22 May 2015 20:53
 Operator : TP/IZ
 Sample : SSTDICC010
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDICC010

Quant Time: May 23 04:47:53 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SIMPAH-BM052215.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat May 23 04:19:45 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	12795	5.00	ng	0.00
6) Naphthalene-d8	10.18	136	47174	5.00	ng	0.00
12) Acenaphthene-d10	14.08	164	27385	5.00	ng	0.00
18) Phenanthrene-d10	16.82	188	96613	5.00	ng	0.00
25) Chrysene-d12	21.04	240	60307	5.00	ng	0.00
32) Perylene-d12	23.14	264	53081	5.00	ng	0.00

System Monitoring Compounds						
4) 2-Fluorophenol	4.99	112	24955	9.95	ng	0.00
5) Phenol-d6	6.60	99	30630	9.92	ng	0.00
7) Nitrobenzene-d5	8.55	82	29825	11.56	ng	0.00
13) 2,4,6-Tribromophenol	15.59	330	8256	8.75	ng	0.00
14) 2-Fluorobiphenyl	12.69	172	79461	9.64	ng	0.00
27) Terphenyl-d14	19.48	244	77789	9.72	ng	-0.01

Target Compounds				Qvalue		
2) 1,4-Dioxane	2.86	88	9441	7.41	ng	98
3) n-Nitrosodimethylamine	3.18	42	16052	10.99	ng	# 97
8) Nitrobenzene	8.60	77	32095	11.87	ng	97
9) Naphthalene	10.22	128	96084	9.51	ng	98
10) Hexachlorobutadiene	10.52	225	19147	9.70	ng	100
11) 2-Methylnaphthalene	11.85	142	67293	9.41	ng	96
15) Acenaphthylene	13.79	152	121195	10.44	ng	100
16) Acenaphthene	14.14	154	69335	9.53	ng	98
23) Anthracene	16.95	178	178768	12.96	ng	98
24) Fluoranthene	18.90	202	168785	6.11	ng	99
26) Pyrene	19.26	202	176064	9.53	ng	100
28) Benzo(a)anthracene	21.02	228	168223	9.66	ng	98
29) Chrysene	21.07	228	147838	8.92	ng	98
30) Bis(2-ethylhexyl)phthalate	20.96	149	103472	11.98	ng	100
31) Indeno(1,2,3-cd)pyrene	25.19	276	161509	8.97	ng	99
33) Benzo(b)fluoranthene	22.51	252	157331	9.63	ng	98
34) Benzo(k)fluoranthene	22.55	252	151958	10.43	ng	98
35) Benzo(a)pyrene	23.04	252	143220	10.21	ng	97
36) Dibenzo(a,h)anthracene	25.21	278	125880	9.76	ng	97
37) Benzo(g,h,i)perylene	25.83	276	128871	9.31	ng	98

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

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