

Data Path : Z:\HPCHEM1\BNA M\DATA\BM101515\
 Data File : BM002906.D
 Acq On : 15 Oct 2015 09:36
 Operator : UM/IZ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SSTDCCC001

Quant Time: Oct 15 10:10:36 2015
 Quant Method : Z:\HPCHEM1\BNA M\Methods\8270-SIM-BM100815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 09 06:47:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.66	152	97789	5.00	ng	-0.02
7) Naphthalene-d8	11.49	136	400771	5.00	ng	-0.03
13) Acenaphthene-d10	15.23	164	209269	5.00	ng	-0.03
19) Phenanthrene-d10	17.95	188	518314	5.00	ng	0.00
26) Chrysene-d12	22.03	240	400347	5.00	ng	-0.03
35) Perylene-d12	24.61	264	372502	5.00	ng	-0.02

System Monitoring Compounds						
4) 2-Fluorophenol	6.13	112	15539	0.74	ng	-0.02
5) Phenol-d6	7.79	99	21700	0.84	ng	-0.02
8) Nitrobenzene-d5	9.86	82	17491	0.87	ng	-0.02
14) 2,4,6-Tribromophenol	16.71	330	6083	0.88	ng	-0.03
15) 2-Fluorobiphenyl	13.88	172	63059	0.99	ng	-0.01
29) Terphenyl-d14	20.48	244	57302	1.07	ng	-0.02

Target Compounds				Qvalue		
2) 1,4-Dioxane	3.77	88	8037	0.91	ng	97
3) n-Nitrosodimethylamine	4.16	42	6571	0.86	ng	# 97
6) bis(2-Chloroethyl)ether	8.04	93	21222	0.83	ng	100
9) Nitrobenzene	9.91	77	18680	0.87	ng	99
10) Naphthalene	11.54	128	81810	0.92	ng	96
11) Hexachlorobutadiene	11.80	225	13315	0.94	ng	100
12) 2-Methylnaphthalene	13.12	142	55263	0.93	ng	96
16) Acenaphthylene	14.96	152	84868	0.96	ng	99
17) Acenaphthene	15.29	154	57057	0.99	ng	# 100
18) Fluorene	16.27	166	69371	0.99	ng	98
20) 4-Bromophenyl-phenylether	17.13	248	20528	0.83	ng	88
21) Hexachlorobenzene	17.28	284	20558	0.68	ng	# 49
22) Pentachlorophenol	17.64	266	1708	0.80	ng	96
23) Phenanthrene	18.00	178	100713	0.66	ng	99
24) Anthracene	18.07	178	92433	0.72	ng	97
25) Fluoranthene	19.96	202	112391	0.67	ng	100
27) Benzidine	20.12	184	44689	0.86	ng	98
28) Pyrene	20.32	202	115974	1.07	ng	100
30) Benzo(a)anthracene	22.01	228	101757	0.89	ng	# 86
31) 3,3'-Dichlorobenzidine	21.93	252	34946	0.79	ng	# 77
32) Chrysene	22.07	228	102218	0.95	ng	# 89
33) Bis(2-ethylhexyl)phthalate	21.85	149	55718	0.80	ng	# 93
34) Indeno(1,2,3-cd)pyrene	27.31	276	102316	0.79	ng	100
36) Benzo(b)fluoranthene	23.82	252	106251	1.00	ng	99
37) Benzo(k)fluoranthene	23.87	252	90399	0.88	ng	99
38) Benzo(a)pyrene	24.49	252	87032	0.89	ng	99
39) Dibenzo(a,h)anthracene	27.31	278	81199	0.85	ng	98
40) Benzo(g,h,i)perylene	28.16	276	84401	0.86	ng	99

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

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