

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089664.D
 Acq On : 24 Feb 2015 6:36
 Operator : TP/JJ
 Sample : SSTDCCC001
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Quant Time: Feb 24 07:45:25 2015

Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE022315.M

Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION

QLast Update : Tue Feb 24 03:46:57 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	33969	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	144407	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	65541	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	117883	5.00	ng	0.00
19) Chrysene-d12	23.66	240	92612	5.00	ng	0.00
25) Perylene-d12	25.34	264	70313	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.60	82	8951	1.11	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	19121	1.20	ng	0.00
21) Terphenyl-d14	22.30	244	21668	1.64	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.70	88	5172	1.19	ng	94
3) n-Nitrosodimethylamine	3.37	42	6862	1.17	ng	# 98
6) Naphthalene	12.29	128	34160	1.16	ng	99
7) 2-Methylnaphthalene	13.94	142	20042	1.21	ng	99
10) Acenaphthylene	16.07	152	116696	1.21	ng	100
11) Acenaphthene	16.49	154	18188	1.16	ng	97
12) Fluorene	17.74	166	20389	1.19	ng	99
14) Hexachlorobenzene	18.99	284	3824	1.19	ng	98
15) Pentachlorophenol	19.43	266	370	0.96	ng	97
16) Phenanthrene	19.81	178	30297	1.17	ng	100
17) Anthracene	19.91	178	26549	1.21	ng	100
18) Fluoranthene	21.72	202	26015	1.20	ng	100
20) Pyrene	22.04	202	28340	1.22	ng	100
22) Benzo(a)anthracene	23.64	228	71105	1.10	ng	100
23) Chrysene	23.68	228	102346	1.17	ng	95
24) Indeno(1,2,3-cd)pyrene	26.74	276	37350	0.93	ng	# 91
26) Benzo(b)fluoranthene	24.91	252	6632	0.87	ng	98
27) Benzo(k)fluoranthene	24.94	252	22236	1.16	ng	99
28) Benzo(a)pyrene	25.27	252	8143	0.93	ng	99
29) Dibenzo(a,h)anthracene	26.77	278	6794	0.93	ng	95
30) Benzo(g,h,i)perylene	27.16	276	41926	1.03	ng	99

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

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