

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089653.D
 Acq On : 23 Feb 2015 22:51
 Operator : TP/JJ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Quant Time: Feb 24 03:18:01 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 03:06:59 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	37405	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	159784	5.00	ng	0.00
8) Acenaphthene-d10	16.43	164	74546	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	132551	5.00	ng	0.00
19) Chrysene-d12	23.66	240	109869	5.00	ng	0.00
25) Perylene-d12	25.34	264	85426	5.00	ng	0.00

System Monitoring Compounds

5) Nitrobenzene-d5	10.61	82	17581	2.42	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	35733	1.84	ng	0.00
21) Terphenyl-d14	22.30	244	40311	2.28	ng	0.00

Target Compounds

						Ovalue
2) 1,4-Dioxane	2.69	88	10418	1.68	ng	98
3) n-Nitrosodimethylamine	3.34	42	13780	2.09	ng	100
6) Naphthalene	12.29	128	65686	1.81	ng	100
7) 2-Methylnaphthalene	13.95	142	38410	1.88	ng	97
10) Acenaphthylene	16.06	152	226702	1.84	ng	100
11) Acenaphthene	16.49	154	35525	1.80	ng	99
12) Fluorene	17.74	166	39175	1.90	ng	100
14) Hexachlorobenzene	18.99	284	7177	1.75	ng	98
15) Pentachlorophenol	19.43	266	936	0.92	ng	95
16) Phenanthrene	19.81	178	58414	1.76	ng	100
17) Anthracene	19.91	178	52816	1.97	ng	100
18) Fluoranthene	21.72	202	53415	2.22	ng	100
20) Pyrene	22.03	202	58405	1.69	ng	100
22) Benzo(a)anthracene	23.65	228	163940	2.03	ng	97
23) Chrysene	23.68	228	204427	1.72	ng	96
24) Indeno(1,2,3-cd)pyrene	26.74	276	101775	2.27	ng	# 92
26) Benzo(b)fluoranthene	24.91	252	20575	2.52	ng	98
27) Benzo(k)fluoranthene	24.94	252	49401	2.30	ng	99
28) Benzo(a)pyrene	25.27	252	22550	2.39	ng	98
29) Dibenzo(a,h)anthracene	26.77	278	19172	2.42	ng	97
30) Benzo(g,h,i)perylene	27.16	276	101860	2.08	ng	97

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

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