

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089651.D
 Acq On : 23 Feb 2015 21:33
 Operator : TP/JJ
 Sample : SSTDICC0.8
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Quant Time: Feb 24 03:17:30 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	40765	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	173619	5.00	ng	0.00
8) Acenaphthene-d10	16.41	164	78060	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	140312	5.00	ng	0.00
19) Chrysene-d12	23.65	240	112266	5.00	ng	0.00
25) Perylene-d12	25.34	264	88106	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.60	82	7000	0.89	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	14565	0.72	ng	0.00
21) Terphenyl-d14	22.30	244	11237	0.62	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.71	88	3797	0.56	ng	97
3) n-Nitrosodimethylamine	3.39	42	4935	0.69	ng	# 96
6) Naphthalene	12.29	128	26983	0.68	ng	99
7) 2-Methylnaphthalene	13.95	142	14431	0.65	ng	97
10) Acenaphthylene	16.06	152	85798	0.66	ng	100
11) Acenaphthene	16.49	154	14202	0.69	ng	99
12) Fluorene	17.74	166	15253	0.70	ng	99
14) Hexachlorobenzene	18.99	284	2967	0.68	ng	98
15) Pentachlorophenol	19.44	266	271	0.25	ng	96
16) Phenanthrene	19.81	178	23892	0.68	ng	100
17) Anthracene	19.91	178	19686	0.69	ng	100
18) Fluoranthene	21.72	202	19595	0.77	ng	99
20) Pyrene	22.03	202	21436	0.61	ng	100
22) Benzo(a)anthracene	23.65	228	56408	0.68	ng	99
23) Chrysene	23.68	228	81378	0.67	ng	96
24) Indeno(1,2,3-cd)pyrene	26.74	276	29027	0.63	ng	# 91
26) Benzo(b)fluoranthene	24.91	252	5682	0.67	ng	99
27) Benzo(k)fluoranthene	24.94	252	16880	0.76	ng	99
28) Benzo(a)pyrene	25.27	252	6408	0.66	ng	98
29) Dibenzo(a,h)anthracene	26.77	278	5365	0.66	ng	96
30) Benzo(g,h,i)perylene	27.15	276	32827	0.65	ng	99

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

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