

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089650.D
 Acq On : 23 Feb 2015 20:54
 Operator : TP/JJ
 Sample : SSTDICC0.5
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Quant Time: Feb 24 03:17:16 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	39700	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	168531	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	74059	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	137633	5.00	ng	0.00
19) Chrysene-d12	23.66	240	105880	5.00	ng	0.00
25) Perylene-d12	25.34	264	81611	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.60	82	3490	0.46	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	8623	0.45	ng	0.00
21) Terphenyl-d14	22.30	244	6350	0.37	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.71	88	2420	0.37	ng	99
3) n-Nitrosodimethylamine	3.41	42	2963	0.42	ng	# 97
6) Naphthalene	12.29	128	16864	0.44	ng	99
7) 2-Methylnaphthalene	13.95	142	9393	0.44	ng	96
10) Acenaphthylene	16.06	152	51449	0.42	ng	99
11) Acenaphthene	16.49	154	8751	0.45	ng	98
12) Fluorene	17.74	166	9521	0.46	ng	100
14) Hexachlorobenzene	18.99	284	1843	0.43	ng	98
15) Pentachlorophenol	19.44	266	164	0.16	ng	92
16) Phenanthrene	19.81	178	15101	0.44	ng	99
17) Anthracene	19.91	178	11780	0.42	ng	99
18) Fluoranthene	21.72	202	11827	0.47	ng	99
20) Pyrene	22.04	202	12892	0.39	ng	100
22) Benzo(a)anthracene	23.64	228	32348	0.42	ng	100
23) Chrysene	23.68	228	50772	0.44	ng	95
24) Indeno(1,2,3-cd)pyrene	26.73	276	14500	0.34	ng	# 84
26) Benzo(b)fluoranthene	24.91	252	2850	0.36	ng	99
27) Benzo(k)fluoranthene	24.94	252	8794	0.43	ng	100
28) Benzo(a)pyrene	25.27	252	3303	0.37	ng	99
29) Dibenzo(a,h)anthracene	26.77	278	2661	0.35	ng	90
30) Benzo(g,h,i)perylene	27.15	276	17066	0.36	ng	96

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

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