

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089648.D
 Acq On : 23 Feb 2015 19:37
 Operator : TP/JJ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Quant Time: Feb 24 03:23:07 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	39660	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	168186	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	75206	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	139827	5.00	ng	0.00
19) Chrysene-d12	23.66	240	105039	5.00	ng	0.00
25) Perylene-d12	25.34	264	80220	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.61	82	890	0.12	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	2099	0.11	ng	0.00
21) Terphenyl-d14	22.30	244	1162	0.07	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.73	88	480	0.07	ng	96
3) n-Nitrosodimethylamine	3.53	42	233m	0.03	ng	
6) Naphthalene	12.29	128	3432	0.09	ng	95
7) 2-Methylnaphthalene	13.95	142	1800	0.08	ng	97
10) Acenaphthylene	16.07	152	9573	0.08	ng	98
11) Acenaphthene	16.49	154	1723	0.09	ng	88
12) Fluorene	17.74	166	1824	0.09	ng	99
14) Hexachlorobenzene	18.99	284	377	0.09	ng	90
15) Pentachlorophenol	19.44	266	64	0.06	ng	# 75
16) Phenanthrene	19.81	178	2907m	0.08	ng	
17) Anthracene	19.91	178	2102	0.07	ng	99
18) Fluoranthene	21.72	202	2013	0.08	ng	100
20) Pyrene	22.04	202	2286	0.07	ng	99
22) Benzo(a)anthracene	23.64	228	6301	0.08	ng	96
23) Chrysene	23.68	228	10089	0.09	ng	95
24) Indeno(1,2,3-cd)pyrene	26.74	276	1606	0.04	ng	# 88
26) Benzo(b)fluoranthene	24.91	252	448	0.06	ng	# 80
27) Benzo(k)fluoranthene	24.94	252	1019m	0.05	ng	
28) Benzo(a)pyrene	25.27	252	442	0.05	ng	# 75
29) Dibenzo(a,h)anthracene	26.77	278	277	0.04	ng	# 52
30) Benzo(g,h,i)perylene	27.15	276	2001	0.04	ng	# 71

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

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