

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089649.D
 Acq On : 23 Feb 2015 20:15
 Operator : TP/JJ
 Sample : SSTDICC0.2
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Quant Time: Feb 24 03:36:43 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	37956	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	161362	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	68401	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	125274	5.00	ng	0.00
19) Chrysene-d12	23.66	240	94850	5.00	ng	0.00
25) Perylene-d12	25.34	264	74241	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.61	82	2732	0.37	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	6078	0.34	ng	0.00
21) Terphenyl-d14	22.30	244	2293	0.15	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.72	88	926	0.15	ng	98
3) n-Nitrosodimethylamine	3.47	42	697	0.10	ng	# 1
6) Naphthalene	12.29	128	6676	0.18	ng	97
7) 2-Methylnaphthalene	13.95	142	3454	0.17	ng	98
10) Acenaphthylene	16.07	152	17992	0.16	ng	100
11) Acenaphthene	16.49	154	3242	0.18	ng	95
12) Fluorene	17.74	166	3398	0.18	ng	100
14) Hexachlorobenzene	19.00	284	657	0.17	ng	90
15) Pentachlorophenol	19.44	266	59	0.06	ng	# 83
16) Phenanthrene	19.81	178	5530	0.18	ng	98
17) Anthracene	19.91	178	4027	0.16	ng	99
18) Fluoranthene	21.72	202	4174	0.18	ng	98
20) Pyrene	22.03	202	4523	0.15	ng	100
22) Benzo(a)anthracene	23.64	228	11179	0.16	ng	98
23) Chrysene	23.68	228	18407	0.18	ng	95
24) Indeno(1,2,3-cd)pyrene	26.74	276	3320	0.09	ng	# 89
26) Benzo(b)fluoranthene	24.91	252	866m	0.12	ng	
27) Benzo(k)fluoranthene	24.94	252	2351m	0.13	ng	
28) Benzo(a)pyrene	25.27	252	968	0.12	ng	# 84
29) Dibenzo(a,h)anthracene	26.77	278	559	0.08	ng	# 68
30) Benzo(g,h,i)perylene	27.15	276	5479	0.13	ng	# 85

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

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