

Data Path : Z:\HPCHEM1\BNA E\DATA\BE022315\
 Data File : BE089656.D
 Acq On : 24 Feb 2015 00:47
 Operator : TP/JJ
 Sample : SSTDICV001
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Quant Time: Feb 24 04:03:36 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 03:46:57 2015

Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	37226	5.00	ng	0.00
4) Naphthalene-d8	12.24	136	157929	5.00	ng	0.00
8) Acenaphthene-d10	16.43	164	70264	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	128274	5.00	ng	0.00
19) Chrysene-d12	23.66	240	104344	5.00	ng	0.00
25) Perylene-d12	25.34	264	80098	5.00	ng	0.00
System Monitoring Compounds						
5) Nitrobenzene-d5	10.60	82	8307	0.96	ng	0.00
9) 2-Fluorobiphenyl	14.87	172	17027	0.98	ng	0.00
21) Terphenyl-d14	22.30	244	16175	1.09	ng	0.00
Target Compounds						Ovalue
2) 1,4-Dioxane	2.70	88	4831	1.01	ng	96
3) n-Nitrosodimethylamine	3.37	42	6653	1.04	ng	97
6) Naphthalene	12.29	128	32335	1.00	ng	100
7) 2-Methylnaphthalene	13.94	142	18667	1.03	ng	100
10) Acenaphthylene	16.06	152	107477	1.04	ng	100
11) Acenaphthene	16.49	154	16982	1.01	ng	99
12) Fluorene	17.74	166	19013	1.03	ng	100
14) Hexachlorobenzene	18.99	284	3613	1.04	ng	97
15) Pentachlorophenol	19.44	266	393	0.94	ng	99
16) Phenanthrene	19.81	178	28714	1.02	ng	100
17) Anthracene	19.91	178	24647	1.03	ng	99
18) Fluoranthene	21.72	202	24712	1.05	ng	99
20) Pyrene	22.04	202	27176	1.04	ng	100
22) Benzo(a)anthracene	23.64	228	72212	0.99	ng	100
23) Chrysene	23.68	228	100333	1.02	ng	95
24) Indeno(1,2,3-cd)pyrene	26.74	276	38013	0.88	ng	# 92
26) Benzo(b)fluoranthene	24.91	252	7294	0.85	ng	98
27) Benzo(k)fluoranthene	24.94	252	22156	1.02	ng	99
28) Benzo(a)pyrene	25.27	252	8520	0.88	ng	99
29) Dibenzo(a,h)anthracene	26.77	278	7055	0.88	ng	98
30) Benzo(g,h,i)perylene	27.16	276	41435	0.92	ng	99

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

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