

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE021815\
 Data File : BE089609.D
 Acq On : 18 Feb 2015 00:55
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.1

Quant Time: Feb 18 15:27:16 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE021815.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 18 05:28:48 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.34	152	44210	5.00	ng	0.00
4) Naphthalene-d8	12.25	136	178973	5.00	ng	0.00
8) Acenaphthene-d10	16.42	164	87193	5.00	ng	0.00
13) Phenanthrene-d10	19.77	188	170015	5.00	ng	0.00
19) Chrysene-d12	23.66	240	162241	5.00	ng	0.00
25) Perylene-d12	25.34	264	142470	5.00	ng	0.00

System Monitoring Compounds						
5) Nitrobenzene-d5	10.61	82	716	0.09	ng	0.00
9) 2-Fluorobiphenyl	14.88	172	2455	0.09	ng	0.00
21) Terphenyl-d14	22.30	244	2401	0.08	ng	0.00

Target Compounds						Qvalue
2) 1,4-Dioxane	2.73	88	556	0.10	ng	97
3) n-Nitrosodimethylamine	3.50	42	373	0.07	ng	# 77
6) Naphthalene	12.30	128	4325	0.10	ng	97
7) 2-Methylnaphthalene	13.95	142	2468	0.09	ng	99
10) Acenaphthylene	16.07	152	13413	0.08	ng	100
11) Acenaphthene	16.50	154	2334	0.09	ng	90
12) Fluorene	17.74	166	2600	0.09	ng	99
14) Hexachlorobenzene	19.00	284	774	0.09	ng	92
16) Phenanthrene	19.82	178	4426	0.09	ng	98
17) Anthracene	19.91	178	3214	0.08	ng	99
18) Fluoranthene	21.72	202	3510	0.08	ng	100
20) Pyrene	22.04	202	3855	0.08	ng	100
22) Benzo(a)anthracene	23.64	228	14049	0.09	ng	97
23) Chrysene	23.69	228	17284	0.09	ng	97
24) Indeno(1,2,3-cd)pyrene	26.75	276	7243	0.05	ng	96
26) Benzo(b)fluoranthene	24.91	252	1531	0.05	ng	# 91
27) Benzo(k)fluoranthene	24.94	252	3204	0.07	ng	# 92
28) Benzo(a)pyrene	25.27	252	1528	0.06	ng	# 89
29) Dibenzo(a,h)anthracene	26.77	278	1429	0.05	ng	# 80
30) Benzo(g,h,i)perylene	27.17	276	8395	0.06	ng	# 88

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

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