

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020615\
 Data File : BE089525.D
 Acq On : 6 Feb 2015 13:10
 Operator : TP/IZ
 Sample : SSTDICC0.2
 Misc :
 ALS Vial : 3 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.2

Manual Integrations
 APPROVED

apatel
 2/9/2015 11:54:48 AM

Quant Time: Feb 07 04:14:18 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SIMPAH-BE020615.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Feb 07 03:25:41 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.43	152	29957	5.00	ng	0.00
2) Naphthalene-d8	12.33	136	111965	5.00	ng	0.00
6) Acenaphthene-d10	16.51	164	41411	5.00	ng	0.00
11) Phenanthrene-d10	19.85	188	66809	5.00	ng	0.00
16) Chrysene-d12	23.72	240	74227	5.00	ng	0.00
22) Perylene-d12	25.42	264	88784	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.69	82	1468	0.19	ng	0.00
7) 2-Fluorobiphenyl	14.97	172	2365	0.18	ng	0.00
18) Terphenyl-d14	22.36	244	1964	0.15	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	12.38	128	5348	0.19	ng	98
5) 2-Methylnaphthalene	14.04	142	2897	0.17	ng	97
8) Acenaphthylene	16.16	152	13297	0.15	ng	99
9) Acenaphthene	16.59	154	2318	0.17	ng	91
10) Fluorene	17.84	166	2289	0.15	ng	99
13) Phenanthrene	19.89	178	3465	0.17	ng	98
14) Anthracene	19.99	178	2551	0.14	ng	97
15) Fluoranthene	21.79	202	2870	0.15	ng	100
17) Pyrene	22.11	202	3015	0.13	ng	98
19) Benzo(a)anthracene	23.71	228	13740	0.17	ng	98
20) Chrysene	23.76	228	14100	0.17	ng	98
21) Indeno(1,2,3-cd)pyrene	26.86	276	20906	0.22	ng	95
23) Benzo(b)fluoranthene	24.98	252	3315	0.17	ng	# 93
24) Benzo(k)fluoranthene	25.01	252	4082	0.14	ng	96
25) Benzo(a)pyrene	25.35	252	3264	0.16	ng	# 89
26) Dibenzo(a,h)anthracene	26.89	278	4260	0.19	ng	# 88
27) Benzo(g,h,i)perylene	27.29	276	18393	0.19	ng	96

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

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