

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE020615\
 Data File : BE089524.D
 Acq On : 6 Feb 2015 12:32
 Operator : TP/IZ
 Sample : SSTDICC0.1
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC0.1

Manual Integrations
 APPROVED

apatel
 2/9/2015 11:54:45 AM

Quant Time: Feb 07 04:13:26 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	32311	5.00	ng	0.00
2) Naphthalene-d8	12.33	136	122716	5.00	ng	0.00
6) Acenaphthene-d10	16.52	164	47715	5.00	ng	0.01
11) Phenanthrene-d10	19.85	188	81375	5.00	ng	0.00
16) Chrysene-d12	23.72	240	90618	5.00	ng	0.00
22) Perylene-d12	25.42	264	104587	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.69	82	790	0.09	ng	0.00
7) 2-Fluorobiphenyl	14.97	172	1342	0.09	ng	0.00
18) Terphenyl-d14	22.36	244	1199	0.07	ng	0.00
Target Compounds						Qvalue
4) Naphthalene	12.38	128	2848	0.09	ng	# 94
5) 2-Methylnaphthalene	14.04	142	1571m	0.08	ng	
8) Acenaphthylene	16.17	152	7424	0.07	ng	# 94
9) Acenaphthene	16.59	154	1318	0.09	ng	82
10) Fluorene	17.84	166	1351	0.08	ng	# 96
13) Phenanthrene	19.90	178	2143	0.09	ng	96
14) Anthracene	19.99	178	1538	0.07	ng	97
15) Fluoranthene	21.79	202	1807	0.08	ng	96
17) Pyrene	22.11	202	1898	0.07	ng	96
19) Benzo(a)anthracene	23.71	228	8598	0.09	ng	96
20) Chrysene	23.75	228	8328	0.08	ng	95
21) Indeno(1,2,3-cd)pyrene	26.85	276	11467	0.10	ng	96
23) Benzo(b)fluoranthene	24.98	252	1933	0.08	ng	# 87
24) Benzo(k)fluoranthene	25.01	252	2401	0.07	ng	# 88
25) Benzo(a)pyrene	25.35	252	1893	0.08	ng	# 82
26) Dibenzo(a,h)anthracene	26.89	278	2240	0.08	ng	# 78
27) Benzo(g,h,i)perylene	27.29	276	10260	0.09	ng	# 88

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

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