

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020515\
 Data File : BE089502.D
 Acq On : 5 Feb 2015 12:39
 Operator : TP/IZ
 Sample : SSTDICC002
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 LabSampleId :
 SSTDICC002

Quant Time: Feb 06 02:48:45 2015
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SIMPAH-BE020515.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 06 02:40:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	9.49	152	26390	5.00	ng	0.00
2) Naphthalene-d8	12.40	136	112274	5.00	ng	0.00
6) Acenaphthene-d10	16.59	164	55379	5.00	ng	0.00
11) Phenanthrene-d10	19.90	188	110897	5.00	ng	0.00
16) Chrysene-d12	23.76	240	121574	5.00	ng	0.00
22) Perylene-d12	25.46	264	119864	5.00	ng	0.00
System Monitoring Compounds						
3) Nitrobenzene-d5	10.76	82	18859	2.20	ng	0.00
7) 2-Fluorobiphenyl	15.03	172	40205	2.30	ng	0.00
18) Terphenyl-d14	22.41	244	51134	2.42	ng	0.00
Target Compounds						Ovalue
4) Naphthalene	12.45	128	60808	2.15	ng	100
5) 2-Methylnaphthalene	14.11	142	43774	2.36	ng	100
8) Acenaphthylene	16.24	152	290044	2.44	ng	100
9) Acenaphthene	16.65	154	43231	2.37	ng	99
10) Fluorene	17.91	166	50774	2.50	ng	98
12) Pentachlorophenol	19.58	266	1359	1.25	ng	96
13) Phenanthrene	19.94	178	79428	2.36	ng	99
14) Anthracene	20.04	178	74513	2.47	ng	100
15) Fluoranthene	21.83	202	79853	2.48	ng	100
17) Pyrene	22.15	202	84735	2.44	ng	100
19) Benzo(a)anthracene	23.75	228	325801	2.44	ng	99
20) Chrysene	23.79	228	326418	2.38	ng	98
21) Indeno(1,2,3-cd)pyrene	26.93	276	375082	2.63	ng	99
23) Benzo(b)fluoranthene	25.02	252	75817	2.72	ng	99
24) Benzo(k)fluoranthene	25.05	252	92794	2.43	ng	98
25) Benzo(a)pyrene	25.39	252	72350	2.56	ng	99
26) Dibenzo(a,h)anthracene	26.95	278	80158	2.66	ng	99
27) Benzo(a,h,i)perylene	27.36	276	309783	2.52	ng	100

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

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