

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE071415\
 Data File : BE090098.D
 Acq On : 14 Jul 2015 14:54
 Operator : TP/IZ
 Sample : SST0.101
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.101

Quant Time: Jul 15 01:06:48 2015
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM02.2-EPA-SIM-BE071415.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jul 15 00:44:52 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1444	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	6492	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.22	164	3685	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	8546	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	7829	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	5826	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	1042	0.10	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	2390	0.10	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.42	128	1838	0.10	ng/ul#	93
5) 2-Methylnaphthalene	12.03	142	1186	0.09	ng/ul	99
7) Acenaphthylene	13.93	152	8029	0.24	ng/ul	95
8) Acenaphthene	14.28	153	5662	0.25	ng/ul	99
9) Fluorene	15.27	166	1670	0.10	ng/ul#	97
12) Phenanthrene	16.99	178	10849	0.24	ng/ul	96
13) Anthracene	17.09	178	10142	0.24	ng/ul	93
15) Fluoranthene	19.00	202	11293	0.23	ng/ul	95
17) Pyrene	19.37	202	11744	0.23	ng/ul	97
18) Benzo(a)anthracene	21.11	228	2184	0.08	ng/ul	96
19) Chrysene	21.16	228	2840	0.10	ng/ul	97
21) Benzo(b)fluoranthene	22.73	252	1665	0.07	ng/ul	88
22) Benzo(k)fluoranthene	22.78	252	2477	0.11	ng/ul#	88
23) Benzo(a)pyrene	23.33	252	1747	0.08	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	25.84	276	7902	0.23	ng/ul	97
25) Dibenzo(a,h)anthracene	25.87	278	1485	0.08	ng/ul	91
26) Benzo(g,h,i)perylene	26.57	276	7227	0.24	ng/ul#	87

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

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