

Data Path : Z:\HPCHEM1\BNA_M\DATA\BM030915\
 Data File : BM000665.D
 Acq On : 09 Mar 2015 17:45
 Operator : TP/IZ
 Sample : SST0.850
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SST0.850

Quant Time: Mar 10 02:36:48 2015
 Quant Method : Z:\HPCHEM1\BNA_M\METHODS\SOM02.2-EPA-SIM-BM030915.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 10 02:25:17 2015
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.67	152	4446	0.40	ng/ul	0.00
2) Naphthalene-d8	10.44	136	14961	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.31	164	4262	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.05	188	14158	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	12662	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	11905	0.40	ng/ul	0.00

System Monitoring Compounds	R.T.	QIon	Response	Conc	Units	Dev(Min)
4) 2-Methylnaphthalene-d10	12.03	152	14164	0.71	ng/ul	-0.01
14) Fluoranthene-d10	19.09	212	23181	0.76	ng/ul	0.00

Target Compounds	R.T.	QIon	Response	Conc	Units	Qvalue
3) Naphthalene	10.49	128	29634	0.78	ng/ul	99
5) 2-Methylnaphthalene	12.11	142	18793	0.71	ng/ul	99
7) Acenaphthylene	14.03	152	23506	1.17	ng/ul	99
8) Acenaphthene	14.36	153	17626	1.15	ng/ul	99
9) Fluorene	15.36	166	21059	1.12	ng/ul	100
11) Pentachlorophenol	16.71	266	810	0.23	ng/ul	96
12) Phenanthrene	17.10	178	31336	0.77	ng/ul	99
13) Anthracene	17.19	178	28142	0.74	ng/ul	98
15) Fluoranthene	19.11	202	34810	0.76	ng/ul	99
17) Pyrene	19.48	202	36268	0.64	ng/ul	99
18) Benzo(a)anthracene	21.23	228	29224	0.73	ng/ul	99
19) Chrysene	21.28	228	34633	0.79	ng/ul	99
21) Benzo(b)fluoranthene	22.79	252	33805	0.76	ng/ul	96
22) Benzo(k)fluoranthene	22.83	252	33587	0.76	ng/ul	96
23) Benzo(a)pyrene	23.35	252	31461	0.78	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.67	276	36462	0.92	ng/ul	100
25) Dibenzo(a,h)anthracene	25.68	278	29610	0.94	ng/ul	96
26) Benzo(g,h,i)perylene	26.33	276	32261	0.93	ng/ul	97

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

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