

Data Path : Z:\HPCHEM1\BNA E\DATA\BE060717\
 Data File : BE093068.D
 Acq On : 7 Jun 2017 14:40
 Operator : SJ/MA
 Sample : SSTD0.855
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.855

Quant Time: Jun 07 15:54:02 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jun 07 14:48:46 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	4116	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	19014	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	10312	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	28379	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	34625	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	26920	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	22738	0.82	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	63043	0.84	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	10.57	128	37851	0.75	ng/ul#	91
5) 2-Methylnaphthalene	12.19	142	26524	0.75	ng/ul	98
7) Acenaphthylene	14.09	152	37368	0.73	ng/ul	96
8) Acenaphthene	14.43	153	29048	0.75	ng/ul	98
9) Fluorene	15.42	166	34104	0.73	ng/ul#	92
11) Pentachlorophenol	16.76	266	3388	0.31	ng/ul	98
12) Phenanthrene	17.14	178	62747	0.69	ng/ul#	86
13) Anthracene	17.22	178	56647	0.66	ng/ul#	86
15) Fluoranthene	19.15	202	78561	0.69	ng/ul	84
17) Pyrene	19.51	202	80264	0.54	ng/ul#	86
18) Benzo(a)anthracene	21.24	228	75352	0.54	ng/ul#	85
19) Chrysene	21.30	228	76916	0.57	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	68584	0.58	ng/ul	86
22) Benzo(k)fluoranthene	22.98	252	71363	0.64	ng/ul#	87
23) Benzo(a)pyrene	23.55	252	64710	0.58	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.18	276	74445	0.51	ng/ul#	92
25) Dibenzo(a,h)anthracene	26.21	278	62555	0.54	ng/ul#	91
26) Benzo(g,h,i)perylene	26.96	276	63664	0.50	ng/ul	97

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

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