

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE061417\
 Data File : BE093178.D
 Acq On : 14 Jun 2017 20:24
 Operator : SJ/MA
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.405

Quant Time: Jun 15 02:29:07 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 15 02:27:21 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.74	152	3073	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	15099	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.37	164	9157	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	23707	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	25132	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	21048	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.11	152	9694	0.44	ng/ul	0.00
14) Fluoranthene-d10	19.13	212	27441	0.43	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.57	128	15383	0.411	ng/ul#	92
5) 2-Methylnaphthalene	12.19	142	11212	0.438	ng/ul	98
7) Acenaphthylene	14.09	152	17629	0.449	ng/ul	96
8) Acenaphthene	14.43	153	13048	0.412	ng/ul	99
9) Fluorene	15.42	166	15205	0.409	ng/ul	93
11) Pentachlorophenol	16.75	266	1929	0.548	ng/ul	97
12) Phenanthrene	17.14	178	26858	0.412	ng/ul#	88
13) Anthracene	17.22	178	25721	0.453	ng/ul#	87
15) Fluoranthene	19.16	202	32597	0.408	ng/ul	83
17) Pyrene	19.52	202	34685	0.468	ng/ul#	82
18) Benzo(a)anthracene	21.25	228	30546	0.456	ng/ul#	88
19) Chrysene	21.30	228	28676	0.412	ng/ul	98
21) Benzo(b)fluoranthene	22.93	252	26408	0.413	ng/ul	94
22) Benzo(k)fluoranthene	22.98	252	26902	0.397	ng/ul	96
23) Benzo(a)pyrene	23.55	252	26013	0.431	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.18	276	28738	0.411	ng/ul#	91
25) Dibenzo(a,h)anthracene	26.20	278	24053	0.408	ng/ul#	90
26) Benzo(g,h,i)perylene	26.95	276	23884	0.398	ng/ul	95

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

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