

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE121817\
 Data File : BE095007.D
 Acq On : 18 Dec 2017 13:26
 Operator : SJ/JU
 Sample : SSTD0.804
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.804

Quant Time: Dec 18 15:33:53 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE121817.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 18 15:33:13 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4688	0.40	ng/ul	0.00
2) Naphthalene-d8	11.29	136	14640	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.05	164	8204	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	18812	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	20654	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	16697	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.85	152	19553	0.83	ng/ul	0.00
14) Fluoranthene-d10	19.73	212	50235	0.78	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.34	128	30899	0.886	ng/ul#	87
5) 2-Methylnaphthalene	12.91	142	22051	0.909	ng/ul	98
7) Acenaphthylene	14.77	152	29945	0.797	ng/ul	94
8) Acenaphthene	15.11	153	22970	0.869	ng/ul	95
9) Fluorene	16.07	166	25303	0.789	ng/ul	90
11) Pentachlorophenol	17.39	266	1418	1.430	ng/ul#	56
12) Phenanthrene	17.78	178	41066	0.813	ng/ul#	81
13) Anthracene	17.87	178	38335	0.745	ng/ul#	86
15) Fluoranthene	19.76	202	55888	0.843	ng/ul	83
17) Pyrene	20.11	202	55234	0.855	ng/ul#	78
18) Benzo(a)anthracene	21.82	228	47354	0.754	ng/ul#	84
19) Chrysene	21.88	228	49551	0.790	ng/ul	95
21) Benzo(b)fluoranthene	23.68	252	47119	0.947	ng/ul	82
22) Benzo(k)fluoranthene	23.73	252	41871	0.791	ng/ul#	84
23) Benzo(a)pyrene	24.38	252	40302	0.815	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.34	276	44065	0.817	ng/ul#	82
25) Dibenzo(a,h)anthracene	27.36	278	36213	0.797	ng/ul#	87
26) Benzo(g,h,i)perylene	28.22	276	37837	0.838	ng/ul	94

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

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