

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE102517\
 Data File : BE094517.D
 Acq On : 24 Oct 2017 16:36
 Operator : SJ/JU
 Sample : SSTD0.147
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.147

Quant Time: Oct 24 19:01:32 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE102517.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Oct 24 18:46:10 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.92	152	1134	0.40	ng/ul	0.00
2) Naphthalene-d8	10.73	136	6200	0.40	ng/ul	0.03
6) Acenaphthene-d10	14.53	164	3816	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	8313	0.40	ng/ul	0.02
16) Chrysene-d12	21.43	240	8715	0.40	ng/ul	0.00
20) Perylene-d12	23.96	264	8457	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.32	152	1020	0.11	ng/ul	0.03
14) Fluoranthene-d10	19.30	212	2456	0.09	ng/ul	0.00

Target Compounds					Qvalue	
3) Naphthalene	10.78	128	1565	0.103	ng/ul#	81
5) 2-Methylnaphthalene	12.40	142	1045	0.106	ng/ul	97
7) Acenaphthylene	14.26	152	1677	0.113	ng/ul	93
8) Acenaphthene	14.60	153	1228	0.100	ng/ul	97
9) Fluorene	15.60	166	1315	0.092	ng/ul	92
12) Phenanthrene	17.31	178	2221	0.102	ng/ul#	91
13) Anthracene	17.41	178	2122	0.100	ng/ul#	94
15) Fluoranthene	19.33	202	2536	0.089	ng/ul	86
17) Pyrene	19.69	202	2731	0.113	ng/ul#	87
18) Benzo(a)anthracene	21.41	228	2491	0.108	ng/ul#	87
19) Chrysene	21.47	228	2664	0.100	ng/ul	95
21) Benzo(b)fluoranthene	23.18	252	2407	0.105	ng/ul	96
22) Benzo(k)fluoranthene	23.23	252	2580	0.097	ng/ul#	95
23) Benzo(a)pyrene	23.85	252	2384	0.100	ng/ul#	95
24) Indeno(1,2,3-cd)pyrene	26.65	276	2508	0.103	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.65	278	2075	0.101	ng/ul#	86
26) Benzo(g,h,i)perylene	27.48	276	2050	0.097	ng/ul#	79

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

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