

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042018\
 Data File : BE096038.D
 Acq On : 20 Apr 2018 9:03
 Operator : SJ/JU
 Sample : SSTDC000.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.443

Quant Time: Apr 21 00:57:08 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Apr 18 01:18:08 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1801	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4596	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2582	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6504	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	8370	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6837	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3205	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9637	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	4710	0.367	ng/ul#	88
5) 2-Methylnaphthalene	12.83	142	3305	0.380	ng/ul	98
7) Acenaphthylene	14.70	152	4962	0.371	ng/ul	97
8) Acenaphthene	15.03	153	3605	0.375	ng/ul	94
9) Fluorene	16.00	166	4087	0.377	ng/ul#	94
11) Pentachlorophenol	17.34	266	364	0.264	ng/ul	94
12) Phenanthrene	17.72	178	7015	0.366	ng/ul#	89
13) Anthracene	17.81	178	6844	0.359	ng/ul#	92
15) Fluoranthene	19.69	202	9391	0.360	ng/ul	85
17) Pyrene	20.04	202	10037	0.353	ng/ul#	81
18) Benzo(a)anthracene	21.75	228	12034	0.356	ng/ul#	85
19) Chrysene	21.81	228	8493	0.349	ng/ul	95
21) Benzo(b)fluoranthene	23.57	252	8039	0.363	ng/ul	88
22) Benzo(k)fluoranthene	23.62	252	8149	0.370	ng/ul#	89
23) Benzo(a)pyrene	24.26	252	7700	0.361	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	27.18	276	8545	0.360	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.19	278	6754	0.354	ng/ul	92
26) Benzo(g,h,i)perylene	28.04	276	7330	0.370	ng/ul	94

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

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