

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041618\
 Data File : BE096000.D
 Acq On : 16 Apr 2018 9:22
 Operator : SJ/JU
 Sample : SSTDC000.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.438

Quant Time: Apr 16 13:01:12 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 10 03:47:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1873	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4853	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2611	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6267	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	9314	0.40	ng/ul	0.00
20) Perylene-d12	24.39	264	7400	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3313	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9620	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	4918	0.363	ng/ul#	90
5) 2-Methylnaphthalene	12.83	142	3288	0.358	ng/ul	98
7) Acenaphthylene	14.70	152	4929	0.364	ng/ul	97
8) Acenaphthene	15.03	153	3602	0.371	ng/ul	95
9) Fluorene	16.00	166	3918	0.357	ng/ul	93
11) Pentachlorophenol	17.34	266	356	0.268	ng/ul	92
12) Phenanthrene	17.72	178	6834	0.370	ng/ul#	88
13) Anthracene	17.81	178	6651	0.362	ng/ul#	92
15) Fluoranthene	19.69	202	9440	0.376	ng/ul	85
17) Pyrene	20.04	202	10865	0.343	ng/ul#	79
18) Benzo(a)anthracene	21.75	228	12937	0.344	ng/ul#	85
19) Chrysene	21.81	228	9134	0.337	ng/ul	97
21) Benzo(b)fluoranthene	23.57	252	8868	0.370	ng/ul	86
22) Benzo(k)fluoranthene	23.63	252	8555	0.359	ng/ul#	86
23) Benzo(a)pyrene	24.27	252	8272	0.358	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.18	276	9217	0.359	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.20	278	7285	0.352	ng/ul	91
26) Benzo(g,h,i)perylene	28.04	276	7864	0.367	ng/ul	95

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

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