

Data Path : Z:\HPCHEM1\BNA E\DATA\BE041618\
 Data File : BE096018.D
 Acq On : 17 Apr 2018 3:14
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.440

Quant Time: Apr 17 05:29:09 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 17 01:14:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1975	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	5262	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	3146	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	7469	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	9574	0.40	ng/ul	0.00
20) Perylene-d12	24.39	264	7487	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	3729	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	10957	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	5375	0.366	ng/ul#	89
5) 2-Methylnaphthalene	12.84	142	3775	0.379	ng/ul	98
7) Acenaphthylene	14.70	152	5948	0.365	ng/ul	97
8) Acenaphthene	15.04	153	4301	0.368	ng/ul	97
9) Fluorene	16.00	166	4843	0.366	ng/ul#	90
11) Pentachlorophenol	17.34	266	631	0.399	ng/ul	95
12) Phenanthrene	17.72	178	8084	0.367	ng/ul#	88
13) Anthracene	17.81	178	8120	0.371	ng/ul#	91
15) Fluoranthene	19.69	202	10572	0.353	ng/ul	83
17) Pyrene	20.04	202	11695	0.359	ng/ul#	79
18) Benzo(a)anthracene	21.75	228	13019	0.337	ng/ul#	84
19) Chrysene	21.81	228	9239	0.332	ng/ul	96
21) Benzo(b)fluoranthene	23.57	252	8934	0.369	ng/ul	91
22) Benzo(k)fluoranthene	23.63	252	8670	0.359	ng/ul	93
23) Benzo(a)pyrene	24.27	252	8351	0.357	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	27.18	276	9603	0.370	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.20	278	7776	0.372	ng/ul	96
26) Benzo(g,h,i)perylene	28.05	276	8083	0.372	ng/ul	95

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

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