

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095465.D
 Acq On : 4 Feb 2018 7:40
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.440

Quant Time: Feb 05 06:41:56 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	1603	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	4665	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	2888	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6898	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6941	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6588	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3821	0.50	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	9613	0.36	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	11.31	128	6319	0.484	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	4419	0.507	ng/ul	99
7) Acenaphthylene	14.74	152	6669	0.480	ng/ul	97
8) Acenaphthene	15.08	153	4877	0.452	ng/ul	96
9) Fluorene	16.03	166	5680	0.511	ng/ul#	86
11) Pentachlorophenol	17.36	266	500	0.374	ng/ul	97
12) Phenanthrene	17.75	178	8892	0.383	ng/ul#	91
13) Anthracene	17.85	178	8312	0.391	ng/ul#	93
15) Fluoranthene	19.72	202	10782	0.359	ng/ul	83
17) Pyrene	20.07	202	11421	0.490	ng/ul#	80
18) Benzo(a)anthracene	21.79	228	10089	0.465	ng/ul#	86
19) Chrysene	21.84	228	9645	0.428	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	9627	0.414	ng/ul	84
22) Benzo(k)fluoranthene	23.67	252	8989	0.423	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	8986	0.433	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	9748	0.428	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.26	278	7820	0.420	ng/ul	92
26) Benzo(g,h,i)perylene	28.11	276	8211	0.426	ng/ul	95

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

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