

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095494.D
 Acq On : 5 Feb 2018 17:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.443

Quant Time: Feb 06 01:32:35 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Feb 06 01:31:24 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1615	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	4256	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	2454	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	5974	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6465	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	6135	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2873	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8170	0.36	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.30	128	4813	0.404	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	3309	0.416	ng/ul	100
7) Acenaphthylene	14.74	152	4942	0.418	ng/ul	98
8) Acenaphthene	15.08	153	3691	0.403	ng/ul	96
9) Fluorene	16.02	166	4279	0.453	ng/ul#	86
11) Pentachlorophenol	17.36	266	383	0.331	ng/ul	97
12) Phenanthrene	17.75	178	7365	0.367	ng/ul#	91
13) Anthracene	17.84	178	6759	0.368	ng/ul#	93
15) Fluoranthene	19.72	202	9260	0.356	ng/ul	85
17) Pyrene	20.07	202	10319	0.476	ng/ul#	79
18) Benzo(a)anthracene	21.78	228	8919	0.441	ng/ul#	84
19) Chrysene	21.84	228	8493	0.405	ng/ul	98
21) Benzo(b)fluoranthene	23.62	252	8375	0.387	ng/ul	86
22) Benzo(k)fluoranthene	23.66	252	7677	0.388	ng/ul#	88
23) Benzo(a)pyrene	24.31	252	7608	0.394	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.23	276	8201	0.387	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.26	278	6607	0.381	ng/ul	94
26) Benzo(g,h,i)perylene	28.11	276	7049	0.393	ng/ul	95

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

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