

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095510.D
 Acq On : 6 Feb 2018 2:44
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.444

Quant Time: Feb 06 03:55:52 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 02:09:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1548	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	4102	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	2447	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6169	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6408	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	6052	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2867	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8134	0.34	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	11.31	128	4654	0.405	ng/ul#	90
5) 2-Methylnaphthalene	12.88	142	3281	0.428	ng/ul	100
7) Acenaphthylene	14.74	152	5027	0.427	ng/ul	98
8) Acenaphthene	15.08	153	3705	0.406	ng/ul	96
9) Fluorene	16.02	166	4355	0.462	ng/ul#	87
11) Pentachlorophenol	17.36	266	455	0.381	ng/ul	93
12) Phenanthrene	17.75	178	7489	0.361	ng/ul#	92
13) Anthracene	17.85	178	6637	0.349	ng/ul#	93
15) Fluoranthene	19.72	202	9228	0.344	ng/ul	85
17) Pyrene	20.07	202	10127	0.471	ng/ul#	81
18) Benzo(a)anthracene	21.78	228	8741	0.436	ng/ul#	83
19) Chrysene	21.84	228	8391	0.403	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	8326	0.390	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	7560	0.387	ng/ul#	88
23) Benzo(a)pyrene	24.31	252	7616	0.399	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	8428	0.403	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.26	278	6834	0.400	ng/ul	91
26) Benzo(g,h,i)perylene	28.11	276	7222	0.408	ng/ul	95

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

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