

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095482.D
 Acq On : 4 Feb 2018 17:49
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.441

Quant Time: Feb 05 06:48:00 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	8.43	152	1501	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	4608	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	2762	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6291	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6660	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6261	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	3674	0.49	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	9077	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	11.31	128	6080	0.471	ng/ul#	90
5) 2-Methylnaphthalene	12.88	142	4258	0.494	ng/ul	99
7) Acenaphthylene	14.74	152	6371	0.479	ng/ul	98
8) Acenaphthene	15.08	153	4646	0.451	ng/ul	96
9) Fluorene	16.02	166	5272	0.496	ng/ul#	85
11) Pentachlorophenol	17.36	266	402	0.330	ng/ul	98
12) Phenanthrene	17.75	178	8648	0.409	ng/ul#	91
13) Anthracene	17.85	178	7799	0.403	ng/ul#	93
15) Fluoranthene	19.72	202	10219	0.373	ng/ul	84
17) Pyrene	20.07	202	10676	0.478	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	9665	0.464	ng/ul#	84
19) Chrysene	21.84	228	9330	0.432	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	9232	0.418	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	8587	0.425	ng/ul#	88
23) Benzo(a)pyrene	24.31	252	8498	0.431	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.25	276	8832	0.408	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.26	278	7100	0.401	ng/ul	93
26) Benzo(g,h,i)perylene	28.12	276	7457	0.407	ng/ul	95

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

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