

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020118\
 Data File : BE095391.D
 Acq On : 2 Feb 2018 11:26
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.433

Quant Time: Feb 02 12:01:13 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jan 27 03:18:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2228	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6818	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	3932	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8572	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9423	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8946	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4696	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	12064	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	7689	0.403	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	5383	0.422	ng/ul	100
7) Acenaphthylene	14.74	152	8140	0.430	ng/ul	96
8) Acenaphthene	15.08	153	5973	0.407	ng/ul	95
9) Fluorene	16.02	166	6491	0.429	ng/ul#	80
11) Pentachlorophenol	17.36	266	731	0.440	ng/ul	94
12) Phenanthrene	17.75	178	11446	0.397	ng/ul#	89
13) Anthracene	17.84	178	11212	0.425	ng/ul#	92
15) Fluoranthene	19.72	202	13321	0.357	ng/ul	83
17) Pyrene	20.07	202	14881	0.471	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	12574	0.427	ng/ul#	86
19) Chrysene	21.84	228	12063	0.394	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	11956	0.379	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	11516	0.399	ng/ul#	86
23) Benzo(a)pyrene	24.32	252	11185	0.397	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	12552	0.406	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.27	278	10178	0.403	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	10612	0.406	ng/ul#	93

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

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