

Data Path : Z:\HPCHEM1\BNA E\DATA\BE101017\
 Data File : BE094272.D
 Acq On : 11 Oct 2017 4:03
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.427

Quant Time: Oct 11 07:35:43 2017
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Oct 11 07:23:41 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1755	0.40	ng/ul	0.00
2) Naphthalene-d8	10.69	136	9024	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.51	164	5020	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	11225	0.40	ng/ul	0.00
16) Chrysene-d12	21.41	240	12117	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	11825	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	5644	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.27	212	14016	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.74	128	8702	0.394	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	5975	0.417	ng/ul	100
7) Acenaphthylene	14.24	152	9234	0.472	ng/ul	96
8) Acenaphthene	14.57	153	6513	0.404	ng/ul	96
9) Fluorene	15.56	166	7405	0.392	ng/ul#	91
11) Pentachlorophenol	16.94	266	362	0.275	ng/ul	90
12) Phenanthrene	17.29	178	11970	0.408	ng/ul#	92
13) Anthracene	17.38	178	12043	0.422	ng/ul#	91
15) Fluoranthene	19.30	202	14294	0.371	ng/ul	84
17) Pyrene	19.66	202	15176	0.450	ng/ul#	82
18) Benzo(a)anthracene	21.39	228	14522	0.452	ng/ul#	85
19) Chrysene	21.45	228	13717	0.370	ng/ul	97
21) Benzo(b)fluoranthene	23.14	252	13977	0.435	ng/ul	92
22) Benzo(k)fluoranthene	23.19	252	13719	0.369	ng/ul	94
23) Benzo(a)pyrene	23.81	252	13393	0.401	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	26.58	276	14907	0.437	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.58	278	12320	0.430	ng/ul	93
26) Benzo(g,h,i)perylene	27.40	276	12037	0.409	ng/ul	96

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

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