

Data Path : U:\HPCHEM1\BNA E\DATA\BE012018\
 Data File : BE095080.D
 Acq On : 20 Jan 2018 9:20
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.408

Quant Time: Jan 22 11:49:29 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE011918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jan 22 04:29:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.46	152	4624	0.40	ng/ul	0.00
2) Naphthalene-d8	11.28	136	14458	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	8104	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	17242	0.40	ng/ul	0.00
16) Chrysene-d12	21.83	240	18279	0.40	ng/ul	0.00
20) Perylene-d12	24.48	264	14702	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.84	152	9592	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	23014	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.33	128	15504	0.405	ng/ul#	89
5) 2-Methylnaphthalene	12.91	142	10704	0.401	ng/ul	99
7) Acenaphthylene	14.77	152	14269	0.393	ng/ul	95
8) Acenaphthene	15.10	153	11280	0.399	ng/ul	93
9) Fluorene	16.05	166	12396	0.411	ng/ul#	80
11) Pentachlorophenol	17.38	266	315	0.309	ng/ul	95
12) Phenanthrene	17.78	178	23001	0.502	ng/ul#	88
13) Anthracene	17.87	178	19224	0.481	ng/ul#	92
15) Fluoranthene	19.75	202	25348	0.433	ng/ul	85
17) Pyrene	20.10	202	26690	0.402	ng/ul#	81
18) Benzo(a)anthracene	21.81	228	19460	0.385	ng/ul#	85
19) Chrysene	21.87	228	22794	0.404	ng/ul	96
21) Benzo(b)fluoranthene	23.66	252	19812	0.400	ng/ul	84
22) Benzo(k)fluoranthene	23.71	252	18398	0.403	ng/ul#	86
23) Benzo(a)pyrene	24.36	252	17050	0.403	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.31	276	15504	0.413	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.34	278	11255	0.420	ng/ul	91
26) Benzo(g,h,i)perylene	28.19	276	15112	0.405	ng/ul	96

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

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