

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095239.D
 Acq On : 29 Jan 2018 10:25
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.419

Quant Time: Jan 29 12:14:17 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 Jan 29 08:11:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2391	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	7373	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4348	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9204	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	10164	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9762	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5044	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13721	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8181	0.396	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	5759	0.418	ng/ul	100
7) Acenaphthylene	14.75	152	8943	0.427	ng/ul	97
8) Acenaphthene	15.08	153	6557	0.404	ng/ul	94
9) Fluorene	16.03	166	6841	0.409	ng/ul#	75
11) Pentachlorophenol	17.36	266	914	0.513	ng/ul	95
12) Phenanthrene	17.75	178	12726	0.411	ng/ul#	86
13) Anthracene	17.85	178	13063	0.461	ng/ul#	91
15) Fluoranthene	19.73	202	15191	0.379	ng/ul	82
17) Pyrene	20.08	202	15304	0.449	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	14112	0.444	ng/ul#	86
19) Chrysene	21.84	228	13369	0.405	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12716	0.369	ng/ul	87
22) Benzo(k)fluoranthene	23.67	252	12673	0.402	ng/ul#	89
23) Benzo(a)pyrene	24.32	252	12463	0.405	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.24	276	13941	0.413	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.27	278	11395	0.413	ng/ul	91
26) Benzo(g,h,i)perylene	28.12	276	11734	0.411	ng/ul	95

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

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