

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031418\
 Data File : BE095756.D
 Acq On : 14 Mar 2018 10:03
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.416

Quant Time: Mar 15 01:28:28 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 14 11:55:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1950	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5751	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3242	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7146	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7207	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	6641	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3552	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8157	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	5857	0.365	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	4066	0.364	ng/ul	99
7) Acenaphthylene	14.73	152	5844	0.357	ng/ul	98
8) Acenaphthene	15.07	153	4368	0.354	ng/ul	93
9) Fluorene	16.03	166	4714	0.352	ng/ul#	89
11) Pentachlorophenol	17.37	266	264	0.263	ng/ul	99
12) Phenanthrene	17.75	178	7644	0.357	ng/ul#	90
13) Anthracene	17.84	178	7101	0.355	ng/ul#	90
15) Fluoranthene	19.72	202	9357	0.361	ng/ul	84
17) Pyrene	20.07	202	9890	0.361	ng/ul#	78
18) Benzo(a)anthracene	21.78	228	8277	0.356	ng/ul#	86
19) Chrysene	21.84	228	8628	0.364	ng/ul	96
21) Benzo(b)fluoranthene	23.61	252	8282	0.360	ng/ul	84
22) Benzo(k)fluoranthene	23.66	252	7706	0.352	ng/ul#	87
23) Benzo(a)pyrene	24.31	252	7522	0.358	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.23	276	8226	0.353	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.25	278	6467	0.347	ng/ul#	90
26) Benzo(g,h,i)perylene	28.11	276	7069	0.353	ng/ul#	94

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

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