

Data Path : U:\HPCHEM1\BNA E\DATA\BE012018\
 Data File : BE095089.D
 Acq On : 20 Jan 2018 14:46
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.409

Quant Time: Jan 22 04:30:53 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	4339	0.40	ng/ul	0.00
2) Naphthalene-d8	11.28	136	14249	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	8129	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	17075	0.40	ng/ul	0.00
16) Chrysene-d12	21.83	240	19290	0.40	ng/ul	0.00
20) Perylene-d12	24.48	264	17585	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.84	152	9568	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	24207	0.46	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.33	128	14849	0.393	ng/ul#	88
5) 2-Methylnaphthalene	12.91	142	10622	0.404	ng/ul	100
7) Acenaphthylene	14.77	152	15268	0.419	ng/ul	96
8) Acenaphthene	15.10	153	11283	0.398	ng/ul	93
9) Fluorene	16.05	166	12363	0.409	ng/ul#	79
11) Pentachlorophenol	17.38	266	883	0.874	ng/ul	94
12) Phenanthrene	17.78	178	22196	0.490	ng/ul#	88
13) Anthracene	17.87	178	21299	0.539	ng/ul#	92
15) Fluoranthene	19.75	202	26091	0.450	ng/ul	84
17) Pyrene	20.10	202	26761	0.382	ng/ul#	80
18) Benzo(a)anthracene	21.81	228	23662	0.443	ng/ul#	85
19) Chrysene	21.87	228	22731	0.382	ng/ul	95
21) Benzo(b)fluoranthene	23.66	252	22965	0.388	ng/ul	83
22) Benzo(k)fluoranthene	23.72	252	20844	0.382	ng/ul#	84
23) Benzo(a)pyrene	24.36	252	21215	0.420	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.31	276	22899	0.510	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.33	278	18480	0.576	ng/ul#	89
26) Benzo(g,h,i)perylene	28.19	276	19536	0.438	ng/ul#	93

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

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