

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095203.D
 Acq On : 28 Jan 2018 12:53
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.416

Quant Time: Jan 29 01:26:15 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 01:25:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2869	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8970	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5438	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	11901	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	13573	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	12851	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	6229	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	17874	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.32	128	10141	0.404	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	7136	0.425	ng/ul	100
7) Acenaphthylene	14.75	152	11200	0.428	ng/ul	95
8) Acenaphthene	15.08	153	8209	0.404	ng/ul	93
9) Fluorene	16.03	166	8934	0.427	ng/ul#	76
11) Pentachlorophenol	17.36	266	1229	0.533	ng/ul	96
12) Phenanthrene	17.75	178	16675	0.417	ng/ul#	88
13) Anthracene	17.85	178	16578	0.453	ng/ul#	92
15) Fluoranthene	19.73	202	20127	0.388	ng/ul	83
17) Pyrene	20.08	202	20155	0.443	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	23428	0.552	ng/ul#	85
19) Chrysene	21.84	228	17931	0.407	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	16863	0.372	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	16723	0.403	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	16195	0.400	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	18191	0.410	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.26	278	14911	0.411	ng/ul#	89
26) Benzo(g,h,i)perylene	28.12	276	15257	0.406	ng/ul	94

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

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