

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012718\
 Data File : BE095196.D
 Acq On : 28 Jan 2018 8:39
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.415

Quant Time: Jan 29 00:17:13 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 00:15:32 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2870	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8985	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	5414	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	12297	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	13621	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	12755	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	6204	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	18256	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	10078	0.401	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	7169	0.427	ng/ul	98
7) Acenaphthylene	14.75	152	11129	0.427	ng/ul	96
8) Acenaphthene	15.08	153	8259	0.409	ng/ul	97
9) Fluorene	16.03	166	8604	0.413	ng/ul#	70
11) Pentachlorophenol	17.36	266	1240	0.521	ng/ul	96
12) Phenanthrene	17.75	178	17339	0.419	ng/ul#	87
13) Anthracene	17.85	178	16819	0.444	ng/ul#	92
15) Fluoranthene	19.73	202	20337	0.380	ng/ul	83
17) Pyrene	20.08	202	22580	0.494	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	18559	0.436	ng/ul#	85
19) Chrysene	21.84	228	17851	0.404	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	16735	0.372	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	16762	0.407	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	16217	0.404	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	18134	0.412	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	14879	0.413	ng/ul#	88
26) Benzo(g,h,i)perylene	28.11	276	15260	0.409	ng/ul#	93

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

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