

Data Path : \\Terastorage\svoasrv\HPCHEM1\BNA_E\Data\BE012618\
 Data File : BE095119.D
 Acq On : 26 Jan 2018 10:04
 Operator : SJ/JU
 Sample : SST00.103
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.103

Quant Time: Jan 26 12:06:51 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 26 12:00:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2901	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8446	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4566	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	10822	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	14974	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13787	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	1330	0.10	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	3999	0.12	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.32	128	2391	0.107	ng/ul#	91
5) 2-Methylnaphthalene	12.89	142	1563	0.100	ng/ul	98
7) Acenaphthylene	14.75	152	2129	0.104	ng/ul	96
8) Acenaphthene	15.08	153	1696	0.107	ng/ul	96
9) Fluorene	16.03	166	1721	0.101	ng/ul#	73
11) Pentachlorophenol	17.36	266	199	0.311	ng/ul	95
12) Phenanthrene	17.75	178	3560	0.124	ng/ul#	86
13) Anthracene	17.85	178	3079	0.123	ng/ul#	91
15) Fluoranthene	19.73	202	4474	0.122	ng/ul	85
17) Pyrene	20.08	202	4994	0.092	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	4555	0.110	ng/ul#	84
19) Chrysene	21.84	228	4716	0.102	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	4729	0.102	ng/ul	91
22) Benzo(k)fluoranthene	23.67	252	4122	0.096	ng/ul	94
23) Benzo(a)pyrene	24.32	252	4108	0.104	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	27.24	276	4451	0.126	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	3581	0.142	ng/ul	94
26) Benzo(g,h,i)perylene	28.11	276	3838	0.110	ng/ul	96

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

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