

Data Path : \\Terastorage\svoasrv\HPCHEM1\BNA_E\Data\BE012618\
 Data File : BE095126.D
 Acq On : 26 Jan 2018 14:12
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.410

Quant Time: Jan 26 14:56:35 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 13:41:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	3038	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8708	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4645	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	11042	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	14979	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13097	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	5606	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	16236	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.32	128	9718	0.399	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	6377	0.392	ng/ul	99
7) Acenaphthylene	14.75	152	8723	0.390	ng/ul	96
8) Acenaphthene	15.08	153	6763	0.390	ng/ul	97
9) Fluorene	16.02	166	7126	0.399	ng/ul#	69
11) Pentachlorophenol	17.36	266	653	0.305	ng/ul	96
12) Phenanthrene	17.75	178	14414	0.388	ng/ul#	85
13) Anthracene	17.84	178	13114	0.386	ng/ul#	90
15) Fluoranthene	19.73	202	18659	0.388	ng/ul	83
17) Pyrene	20.08	202	19880	0.396	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	18146	0.387	ng/ul#	85
19) Chrysene	21.84	228	19661	0.404	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	18734	0.405	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	16841	0.398	ng/ul#	86
23) Benzo(a)pyrene	24.32	252	16295	0.395	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.25	276	17982	0.397	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	14637	0.395	ng/ul#	87
26) Benzo(g,h,i)perylene	28.12	276	15242	0.398	ng/ul#	93

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

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