

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095241.D
 Acq On : 29 Jan 2018 11:43
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.420

Quant Time: Jan 30 04:39:47 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 08:11:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2400	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7232	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4482	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9386	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10381	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9919	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5058	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13745	0.38	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.31	128	8162	0.403	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5754	0.426	ng/ul	99
7) Acenaphthylene	14.75	152	9040	0.419	ng/ul	97
8) Acenaphthene	15.08	153	6610	0.395	ng/ul	93
9) Fluorene	16.03	166	6881	0.399	ng/ul#	75
11) Pentachlorophenol	17.36	266	829	0.456	ng/ul	97
12) Phenanthrene	17.75	178	13401	0.425	ng/ul#	88
13) Anthracene	17.85	178	13022	0.451	ng/ul#	92
15) Fluoranthene	19.73	202	15191	0.372	ng/ul	81
17) Pyrene	20.08	202	15198	0.436	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	14251	0.439	ng/ul#	86
19) Chrysene	21.84	228	13542	0.402	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	12939	0.370	ng/ul	86
22) Benzo(k)fluoranthene	23.67	252	12782	0.399	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	12662	0.405	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.25	276	14123	0.412	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	11472	0.409	ng/ul#	90
26) Benzo(g,h,i)perylene	28.12	276	11806	0.407	ng/ul	94

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

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