

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095256.D
 Acq On : 29 Jan 2018 20:43
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.421

Quant Time: Jan 30 04:45:58 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jan 30 04:43:36 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2323	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6943	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4170	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8875	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9927	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9241	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4871	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12828	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.31	128	7747	0.398	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5521	0.425	ng/ul	99
7) Acenaphthylene	14.75	152	8685	0.433	ng/ul	96
8) Acenaphthene	15.08	153	6319	0.406	ng/ul	95
9) Fluorene	16.02	166	6746	0.420	ng/ul#	76
11) Pentachlorophenol	17.36	266	800	0.465	ng/ul	96
12) Phenanthrene	17.75	178	12455	0.417	ng/ul#	87
13) Anthracene	17.84	178	12313	0.451	ng/ul#	92
15) Fluoranthene	19.72	202	14376	0.372	ng/ul	84
17) Pyrene	20.08	202	15788	0.474	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	13304	0.429	ng/ul#	86
19) Chrysene	21.84	228	12807	0.397	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	12121	0.372	ng/ul	86
22) Benzo(k)fluoranthene	23.67	252	12328	0.413	ng/ul#	86
23) Benzo(a)pyrene	24.32	252	11797	0.405	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.25	276	13186	0.413	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.27	278	10720	0.410	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	11104	0.411	ng/ul#	94

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

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