

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE012918\
 Data File : BE095273.D
 Acq On : 30 Jan 2018 6:52
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.422

Quant Time: Jan 30 07:28:50 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 06:14:31 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2245	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6550	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3783	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8212	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9622	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	8961	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4473	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12297	0.39	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.31	128	7353	0.401	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	5122	0.418	ng/ul	99
7) Acenaphthylene	14.74	152	7683	0.422	ng/ul	97
8) Acenaphthene	15.07	153	5694	0.403	ng/ul	92
9) Fluorene	16.02	166	6102	0.419	ng/ul#	78
11) Pentachlorophenol	17.36	266	727	0.457	ng/ul	94
12) Phenanthrene	17.75	178	11728	0.425	ng/ul#	88
13) Anthracene	17.84	178	11282	0.446	ng/ul#	92
15) Fluoranthene	19.72	202	13845	0.387	ng/ul	84
17) Pyrene	20.08	202	14226	0.441	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	12942	0.430	ng/ul#	86
19) Chrysene	21.84	228	12533	0.401	ng/ul	95
21) Benzo(b)fluoranthene	23.62	252	12091	0.382	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	11695	0.404	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	11626	0.412	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	12852	0.415	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.27	278	10416	0.411	ng/ul#	89
26) Benzo(g,h,i)perylene	28.11	276	10767	0.411	ng/ul	94

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

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