

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095275.D
 Acq On : 30 Jan 2018 8:51
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.423

Quant Time: Jan 31 01:18:52 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	2215	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6776	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4114	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8757	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9646	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9066	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4749	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12322	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	7684	0.405	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5464	0.431	ng/ul	99
7) Acenaphthylene	14.74	152	8224	0.415	ng/ul	97
8) Acenaphthene	15.08	153	6155	0.401	ng/ul	94
9) Fluorene	16.03	166	6597	0.417	ng/ul#	75
11) Pentachlorophenol	17.36	266	672	0.396	ng/ul	97
12) Phenanthrene	17.75	178	12366	0.420	ng/ul#	88
13) Anthracene	17.85	178	11595	0.430	ng/ul#	92
15) Fluoranthene	19.72	202	13921	0.365	ng/ul	84
17) Pyrene	20.08	202	15040	0.465	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	12935	0.429	ng/ul#	86
19) Chrysene	21.84	228	12724	0.406	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12306	0.385	ng/ul	84
22) Benzo(k)fluoranthene	23.67	252	11801	0.403	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	11753	0.412	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	13094	0.418	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.27	278	10552	0.412	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	10951	0.413	ng/ul#	93

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

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