

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013118\
 Data File : BE095322.D
 Acq On : 31 Jan 2018 18:11
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Feb 01 04:01:39 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 08:16:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2416	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7478	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4390	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8975	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9648	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9423	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5221	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12636	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8385	0.400	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5993	0.429	ng/ul	99
7) Acenaphthylene	14.75	152	9085	0.430	ng/ul	96
8) Acenaphthene	15.08	153	6582	0.402	ng/ul	94
9) Fluorene	16.02	166	6958	0.412	ng/ul#	76
11) Pentachlorophenol	17.36	266	765	0.440	ng/ul	96
12) Phenanthrene	17.75	178	12412	0.411	ng/ul#	87
13) Anthracene	17.84	178	11717	0.424	ng/ul#	93
15) Fluoranthene	19.72	202	14169	0.363	ng/ul	84
17) Pyrene	20.08	202	15553	0.480	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	13336	0.442	ng/ul#	85
19) Chrysene	21.84	228	12625	0.403	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12281	0.369	ng/ul	86
22) Benzo(k)fluoranthene	23.67	252	12091	0.398	ng/ul#	88
23) Benzo(a)pyrene	24.32	252	11898	0.401	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	13338	0.410	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.27	278	10814	0.406	ng/ul#	89
26) Benzo(g,h,i)perylene	28.12	276	11182	0.406	ng/ul	95

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

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