

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013118\
 Data File : BE095338.D
 Acq On : 1 Feb 2018 3:43
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Feb 01 06:25:10 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 04:05:22 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2482	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7786	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4492	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9358	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9991	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9598	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5298	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13274	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8648	0.397	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	6076	0.417	ng/ul	99
7) Acenaphthylene	14.75	152	9267	0.429	ng/ul	97
8) Acenaphthene	15.08	153	6803	0.406	ng/ul	95
9) Fluorene	16.02	166	7179	0.415	ng/ul#	78
11) Pentachlorophenol	17.36	266	864	0.477	ng/ul	98
12) Phenanthrene	17.75	178	13028	0.414	ng/ul#	88
13) Anthracene	17.84	178	12607	0.438	ng/ul#	92
15) Fluoranthene	19.72	202	14809	0.363	ng/ul	84
17) Pyrene	20.08	202	16229	0.484	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	13689	0.438	ng/ul#	86
19) Chrysene	21.84	228	13020	0.401	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12797	0.378	ng/ul	84
22) Benzo(k)fluoranthene	23.67	252	12439	0.402	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	12240	0.405	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	13672	0.412	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	11093	0.409	ng/ul#	89
26) Benzo(g,h,i)perylene	28.12	276	11431	0.407	ng/ul	95

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

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