

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013118\
 Data File : BE095355.D
 Acq On : 1 Feb 2018 13:51
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 34 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :

Quant Time: Feb 01 14:26:02 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 06:27:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2334	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7296	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4211	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8786	0.40	ng/ul	0.00
16) Chrysene-d12	21.81	240	9484	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9139	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5044	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12309	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8167	0.400	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5710	0.419	ng/ul	99
7) Acenaphthylene	14.75	152	8763	0.432	ng/ul	97
8) Acenaphthene	15.08	153	6313	0.402	ng/ul	95
9) Fluorene	16.03	166	6835	0.422	ng/ul#	81
11) Pentachlorophenol	17.36	266	741	0.435	ng/ul	96
12) Phenanthrene	17.75	178	11814	0.400	ng/ul#	89
13) Anthracene	17.85	178	11150	0.412	ng/ul#	93
15) Fluoranthene	19.72	202	13676	0.357	ng/ul	84
17) Pyrene	20.08	202	14774	0.464	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	12909	0.435	ng/ul#	86
19) Chrysene	21.84	228	12287	0.399	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12055	0.374	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	11703	0.397	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	11440	0.397	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	12860	0.407	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.27	278	10401	0.403	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	10792	0.404	ng/ul	95

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

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