

Data Path : U:\HPCHEM1\BNA E\DATA\BE012618\
 Data File : BE095122.D
 Acq On : 26 Jan 2018 11:48
 Operator : SJ/JU
 Sample : SST00.806
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST00.806

Quant Time: Jan 26 13:03:19 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 26 12:06:50 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2954	0.40	ng/ul	0.00
2) Naphthalene-d8	11.27	136	8353	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4569	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.73	188	10814	0.40	ng/ul	0.02
16) Chrysene-d12	21.80	240	15155	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	13016	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	11021	0.80	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	33239	0.99	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.32	128	18781	0.849	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	12609	0.817	ng/ul	99
7) Acenaphthylene	14.74	152	17731	0.866	ng/ul	96
8) Acenaphthene	15.08	153	13606	0.855	ng/ul	95
9) Fluorene	16.02	166	14024	0.825	ng/ul#	66
11) Pentachlorophenol	17.36	266	1525	2.384	ng/ul	96
12) Phenanthrene	17.75	178	28058	0.977	ng/ul#	85
13) Anthracene	17.85	178	26816	1.071	ng/ul#	90
15) Fluoranthene	19.73	202	38270	1.043	ng/ul	82
17) Pyrene	20.08	202	39988	0.726	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	37726	0.900	ng/ul#	84
19) Chrysene	21.85	228	39461	0.843	ng/ul	98
21) Benzo(b)fluoranthene	23.62	252	37954	0.866	ng/ul	82
22) Benzo(k)fluoranthene	23.68	252	34208	0.846	ng/ul#	84
23) Benzo(a)pyrene	24.32	252	33436	0.894	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	27.25	276	36886	1.110	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.27	278	30278	1.276	ng/ul#	86
26) Benzo(g,h,i)perylene	28.12	276	30943	0.937	ng/ul#	92

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

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