

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031318\
 Data File : BE095750.D
 Acq On : 13 Mar 2018 15:09
 Operator : SJ/JU
 Sample : SST0.811
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.811

Quant Time: Mar 13 16:25:31 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Mar 13 16:24:15 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1776	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5104	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	2846	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	6365	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6480	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	5677	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	6977	0.84	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	16332	0.67	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.30	128	11379	0.796	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	8041	0.843	ng/ul	99
7) Acenaphthylene	14.73	152	11728	0.856	ng/ul	96
8) Acenaphthene	15.07	153	8761	0.825	ng/ul	94
9) Fluorene	16.04	166	9625	0.879	ng/ul#	93
11) Pentachlorophenol	17.37	266	660	0.535	ng/ul	96
12) Phenanthrene	17.75	178	15291	0.714	ng/ul#	89
13) Anthracene	17.84	178	14571	0.744	ng/ul#	91
15) Fluoranthene	19.72	202	18783	0.678	ng/ul	83
17) Pyrene	20.07	202	19385	0.892	ng/ul#	78
18) Benzo(a)anthracene	21.78	228	16785	0.828	ng/ul#	85
19) Chrysene	21.84	228	17045	0.810	ng/ul	96
21) Benzo(b)fluoranthene	23.61	252	16226	0.810	ng/ul	83
22) Benzo(k)fluoranthene	23.66	252	15444	0.843	ng/ul#	85
23) Benzo(a)pyrene	24.30	252	14871	0.832	ng/ul#	77
24) Indeno(1,2,3-cd)pyrene	27.23	276	16615	0.847	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.25	278	13384	0.834	ng/ul#	86
26) Benzo(g,h,i)perylene	28.10	276	14160	0.853	ng/ul#	91

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

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