

Data Path : Z:\HPCHEM1\BNA E\DATA\BE040618\
 Data File : BE095956.D
 Acq On : 6 Apr 2018 14:51
 Operator : SJ/JU
 Sample : SST0.832
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SST0.832

Quant Time: Apr 06 15:52:12 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 06 15:00:12 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1822	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4659	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.98	164	2634	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6373	0.40	ng/ul	0.00
16) Chrysene-d12	21.78	240	9103	0.40	ng/ul	0.00
20) Perylene-d12	24.39	264	7011	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	6845	0.87	ng/ul	0.00
14) Fluoranthene-d10	19.67	212	20881	1.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.26	128	10214	0.785	ng/ul#	88
5) 2-Methylnaphthalene	12.84	142	7021	0.777	ng/ul	100
7) Acenaphthylene	14.70	152	10901	0.820	ng/ul	96
8) Acenaphthene	15.04	153	7729	0.772	ng/ul	95
9) Fluorene	16.00	166	8712	0.802	ng/ul#	89
11) Pentachlorophenol	17.34	266	1011	1.129	ng/ul	94
12) Phenanthrene	17.73	178	14846	0.778	ng/ul#	88
13) Anthracene	17.82	178	15003	0.841	ng/ul#	91
15) Fluoranthene	19.69	202	20224	0.875	ng/ul	84
17) Pyrene	20.05	202	21781	0.630	ng/ul#	78
18) Benzo(a)anthracene	21.76	228	26197	0.892	ng/ul#	84
19) Chrysene	21.81	228	18991	0.635	ng/ul	96
21) Benzo(b)fluoranthene	23.59	252	17820	0.734	ng/ul	84
22) Benzo(k)fluoranthene	23.64	252	18212	0.789	ng/ul#	87
23) Benzo(a)pyrene	24.28	252	17306	0.781	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.20	276	19533	0.794	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.21	278	16023	0.815	ng/ul#	88
26) Benzo(g,h,i)perylene	28.07	276	16262	0.770	ng/ul#	94

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

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