

Data Path : Z:\HPCHEM1\BNA E\DATA\BE050318\
 Data File : BE096321.D
 Acq On : 3 May 2018 11:20
 Operator : SJ/JU
 Sample : SSTD0.873
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.873

Quant Time: May 03 11:57:45 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE050318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu May 03 11:56:34 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	557	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2134	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	1232	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3427	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	4618	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3611	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.73	152	3114	0.82	ng/ul	0.00
14) Fluoranthene-d10	19.63	212	11612	0.79	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.22	128	4584	0.773	ng/ul#	89
5) 2-Methylnaphthalene	12.81	142	3106	0.807	ng/ul	99
7) Acenaphthylene	14.67	152	5209	0.825	ng/ul	97
8) Acenaphthene	15.00	153	3728	0.812	ng/ul	94
9) Fluorene	15.97	166	5343	0.886	ng/ul	96
11) Pentachlorophenol	17.31	266	341	0.736	ng/ul	95
12) Phenanthrene	17.69	178	8190	0.790	ng/ul#	88
13) Anthracene	17.78	178	8137	0.829	ng/ul#	90
15) Fluoranthene	19.66	202	11673	0.787	ng/ul	83
17) Pyrene	20.02	202	3555	0.639	ng/ul#	76
18) Benzo(a)anthracene	21.73	228	5859	0.788	ng/ul#	87
19) Chrysene	21.78	228	10773	0.781	ng/ul	95
21) Benzo(b)fluoranthene	23.54	252	9916	0.810	ng/ul	85
22) Benzo(k)fluoranthene	23.59	252	9584	0.804	ng/ul#	86
23) Benzo(a)pyrene	24.23	252	9232	0.799	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.12	276	10334	0.814	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.13	278	8440	0.813	ng/ul#	90
26) Benzo(g,h,i)perylene	27.98	276	8762	0.822	ng/ul	95

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

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