

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042018\
 Data File : BE096044.D
 Acq On : 20 Apr 2018 17:01
 Operator : SJ/JU
 Sample : SSTDC000.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.444

Quant Time: Apr 21 01:00:17 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 21 00:59:26 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1872	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4920	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2907	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6882	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	8426	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6635	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3518	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9776	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	5056	0.368	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	3525	0.378	ng/ul	99
7) Acenaphthylene	14.70	152	5408	0.359	ng/ul	98
8) Acenaphthene	15.03	153	3968	0.367	ng/ul	95
9) Fluorene	15.99	166	4509	0.369	ng/ul#	89
11) Pentachlorophenol	17.33	266	498	0.342	ng/ul	91
12) Phenanthrene	17.72	178	7406	0.365	ng/ul#	88
13) Anthracene	17.81	178	7379	0.366	ng/ul#	92
15) Fluoranthene	19.69	202	9434	0.342	ng/ul	83
17) Pyrene	20.04	202	11055	0.386	ng/ul#	80
18) Benzo(a)anthracene	21.75	228	11026	0.324	ng/ul#	86
19) Chrysene	21.81	228	8191	0.334	ng/ul	97
21) Benzo(b)fluoranthene	23.57	252	7876	0.367	ng/ul	90
22) Benzo(k)fluoranthene	23.62	252	7867	0.368	ng/ul	92
23) Benzo(a)pyrene	24.26	252	7573	0.366	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	27.17	276	8489	0.369	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.19	278	6808	0.367	ng/ul	95
26) Benzo(g,h,i)perylene	28.04	276	7115	0.370	ng/ul	96

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

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