

Data Path : Z:\SVOASRV\HPCHEM1\BNA_E\DATA\BE070618\
 Data File : BE096874.D
 Acq On : 7 Jul 2018 4:05
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.452

Quant Time: Jul 07 06:49:15 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE070618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 06 16:55:25 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	2517	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	13700	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.03	164	7852	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.78	188	18786	0.40	ng/ul	0.00
16) Chrysene-d12	20.98	240	20328	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	21548	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.74	152	9259	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	22158	0.47	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.19	128	13239	0.400	ng/ul#	91
5) 2-Methylnaphthalene	11.81	142	9341	0.414	ng/ul	99
7) Acenaphthylene	13.74	152	14157	0.440	ng/ul	97
8) Acenaphthene	14.09	153	10567	0.398	ng/ul	97
9) Fluorene	15.08	166	11285	0.412	ng/ul	92
11) Pentachlorophenol	16.43	266	1449	1.178	ng/ul	96
12) Phenanthrene	16.82	178	19217	0.397	ng/ul#	90
13) Anthracene	16.91	178	19301	0.469	ng/ul#	93
15) Fluoranthene	18.85	202	22915	0.444	ng/ul	84
17) Pyrene	19.21	202	23948	0.359	ng/ul#	84
18) Benzo(a)anthracene	20.96	228	23322	0.486	ng/ul#	83
19) Chrysene	21.02	228	22517	0.362	ng/ul	97
21) Benzo(b)fluoranthene	22.54	252	24844	0.422	ng/ul	88
22) Benzo(k)fluoranthene	22.58	252	23589	0.367	ng/ul#	90
23) Benzo(a)pyrene	23.12	252	23371	0.470	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.50	276	26439	0.512	ng/ul#	81
25) Dibenzo(a,h)anthracene	25.53	278	22225	0.542	ng/ul	91
26) Benzo(g,h,i)perylene	26.20	276	22208	0.443	ng/ul	97

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

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