

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096182.D
 Acq On : 26 Apr 2018 16:54
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.466

Quant Time: Apr 27 02:33:30 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Apr 27 02:29:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	631	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2322	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1352	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3439	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	4437	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3465	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1761	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5419	0.34	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.23	128	2626	0.406	ng/ul#	92
5) 2-Methylnaphthalene	12.81	142	1783	0.412	ng/ul	99
7) Acenaphthylene	14.67	152	2764	0.397	ng/ul	97
8) Acenaphthene	15.01	153	2035	0.405	ng/ul	92
9) Fluorene	15.97	166	2662	0.400	ng/ul	93
11) Pentachlorophenol	17.31	266	206	0.249	ng/ul	92
12) Phenanthrene	17.70	178	4616	0.456	ng/ul#	89
13) Anthracene	17.78	178	4062	0.400	ng/ul#	91
15) Fluoranthene	19.67	202	6438	0.416	ng/ul	84
17) Pyrene	20.02	202	3078	0.346	ng/ul#	78
18) Benzo(a)anthracene	21.73	228	4805	0.358	ng/ul#	86
19) Chrysene	21.79	228	5306	0.430	ng/ul	97
21) Benzo(b)fluoranthene	23.54	252	5140	0.392	ng/ul	86
22) Benzo(k)fluoranthene	23.59	252	4573	0.365	ng/ul#	91
23) Benzo(a)pyrene	24.23	252	4415	0.387	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	27.13	276	4956	0.323	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.14	278	4009	0.312	ng/ul	93
26) Benzo(g,h,i)perylene	28.00	276	4128	0.309	ng/ul	95

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

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