

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096076.D
 Acq On : 23 Apr 2018 9:48
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.447

Quant Time: Apr 24 04:56:16 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Apr 23 01:07:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1642	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	4183	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2409	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	5764	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7352	0.40	ng/ul	0.00
20) Perylene-d12	24.37	264	6192	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	2909	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	8388	0.35	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	4301	0.368	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	3022	0.382	ng/ul	100
7) Acenaphthylene	14.70	152	4592	0.368	ng/ul	98
8) Acenaphthene	15.03	153	3346	0.373	ng/ul	95
9) Fluorene	15.99	166	3734	0.369	ng/ul	90
11) Pentachlorophenol	17.33	266	315	0.258	ng/ul	96
12) Phenanthrene	17.72	178	6127	0.361	ng/ul#	89
13) Anthracene	17.80	178	6120	0.362	ng/ul#	91
15) Fluoranthene	19.68	202	8109	0.351	ng/ul	85
17) Pyrene	20.04	202	8659	0.347	ng/ul#	80
18) Benzo(a)anthracene	21.74	228	9777	0.329	ng/ul#	85
19) Chrysene	21.80	228	7386	0.346	ng/ul	96
21) Benzo(b)fluoranthene	23.57	252	7552	0.377	ng/ul	86
22) Benzo(k)fluoranthene	23.62	252	7068	0.354	ng/ul#	91
23) Benzo(a)pyrene	24.26	252	6972	0.361	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	27.17	276	8003	0.372	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.18	278	6407	0.370	ng/ul#	91
26) Benzo(g,h,i)perylene	28.04	276	6692	0.373	ng/ul	96

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

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