

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040918\
 Data File : BE095966.D
 Acq On : 9 Apr 2018 15:55
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.437

Quant Time: Apr 10 03:48:06 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Apr 10 03:47:21 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1936	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	5000	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2744	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6640	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	8896	0.40	ng/ul	0.00
20) Perylene-d12	24.40	264	7232	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3639	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	10507	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.26	128	5474	0.392	ng/ul#	89
5) 2-Methylnaphthalene	12.83	142	3715	0.392	ng/ul	100
7) Acenaphthylene	14.70	152	5644	0.397	ng/ul	98
8) Acenaphthene	15.04	153	4029	0.395	ng/ul	95
9) Fluorene	16.00	166	4538	0.394	ng/ul	91
11) Pentachlorophenol	17.34	266	418	0.297	ng/ul	95
12) Phenanthrene	17.73	178	7703	0.394	ng/ul#	89
13) Anthracene	17.81	178	7663	0.393	ng/ul#	91
15) Fluoranthene	19.69	202	10332	0.388	ng/ul	85
17) Pyrene	20.04	202	11703	0.387	ng/ul#	81
18) Benzo(a)anthracene	21.75	228	13934	0.388	ng/ul#	84
19) Chrysene	21.82	228	9677	0.374	ng/ul	97
21) Benzo(b)fluoranthene	23.58	252	9726	0.416	ng/ul	85
22) Benzo(k)fluoranthene	23.63	252	8946	0.384	ng/ul#	88
23) Benzo(a)pyrene	24.27	252	8899	0.394	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.19	276	9727	0.388	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.21	278	7935	0.393	ng/ul#	89
26) Benzo(g,h,i)perylene	28.07	276	8217	0.392	ng/ul	95

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

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