

Data Path : U:\HPCHEM1\BNA E\DATA\BE020218\
 Data File : BE095393.D
 Acq On : 2 Feb 2018 12:38
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.434

Quant Time: Feb 02 15:24:42 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Feb 02 04:53:51 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2392	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7536	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4300	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9558	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9657	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9125	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5128	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	12333	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8452	0.401	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	5913	0.420	ng/ul	100
7) Acenaphthylene	14.74	152	8759	0.423	ng/ul	98
8) Acenaphthene	15.08	153	6443	0.401	ng/ul	97
9) Fluorene	16.03	166	7259	0.439	ng/ul#	84
11) Pentachlorophenol	17.36	266	671	0.362	ng/ul	95
12) Phenanthrene	17.75	178	11856	0.369	ng/ul#	91
13) Anthracene	17.85	178	11242	0.382	ng/ul#	92
15) Fluoranthene	19.72	202	13907	0.334	ng/ul	83
17) Pyrene	20.07	202	14992	0.463	ng/ul#	81
18) Benzo(a)anthracene	21.79	228	12922	0.428	ng/ul#	86
19) Chrysene	21.84	228	12406	0.396	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	12400	0.385	ng/ul	84
22) Benzo(k)fluoranthene	23.67	252	11696	0.397	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	11626	0.404	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.24	276	12773	0.405	ng/ul#	76
25) Dibenzo(a,h)anthracene	27.26	278	10242	0.397	ng/ul#	87
26) Benzo(g,h,i)perylene	28.12	276	10786	0.404	ng/ul	95

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

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