

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042318\
 Data File : BE096105.D
 Acq On : 24 Apr 2018 3:13
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.449

Quant Time: Apr 24 16:50:09 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 24 06:59:54 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.38	152	1873	0.40	ng/ul	0.00
2) Naphthalene-d8	11.20	136	5108	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.97	164	2870	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6374	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7487	0.40	ng/ul	0.00
20) Perylene-d12	24.38	264	6526	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.76	152	3520	0.37	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	9044	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.25	128	5292	0.371	ng/ul#	88
5) 2-Methylnaphthalene	12.83	142	3655	0.378	ng/ul	100
7) Acenaphthylene	14.70	152	5531	0.372	ng/ul	97
8) Acenaphthene	15.03	153	3912	0.366	ng/ul	95
9) Fluorene	15.99	166	4284	0.355	ng/ul	86
11) Pentachlorophenol	17.34	266	385	0.285	ng/ul#	90
12) Phenanthrene	17.72	178	6933	0.369	ng/ul#	89
13) Anthracene	17.81	178	7032	0.376	ng/ul#	92
15) Fluoranthene	19.69	202	8757	0.343	ng/ul	84
17) Pyrene	20.04	202	9472	0.372	ng/ul#	80
18) Benzo(a)anthracene	21.75	228	11124	0.368	ng/ul#	86
19) Chrysene	21.81	228	7671	0.352	ng/ul	97
21) Benzo(b)fluoranthene	23.57	252	7851	0.372	ng/ul	89
22) Benzo(k)fluoranthene	23.62	252	7276	0.346	ng/ul	94
23) Benzo(a)pyrene	24.26	252	7300	0.358	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	27.18	276	8591	0.379	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.19	278	6942	0.381	ng/ul	93
26) Benzo(g,h,i)perylene	28.05	276	7050	0.373	ng/ul	95

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

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