

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042718\
 Data File : BE096211.D
 Acq On : 27 Apr 2018 17:48
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.469

Quant Time: Apr 27 23:59:59 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 23:57:47 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.35	152	707	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2691	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.94	164	1501	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3740	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	5244	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	4031	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	1947	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5937	0.34	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	2969	0.396	ng/ul#	91
5) 2-Methylnaphthalene	12.81	142	2020	0.402	ng/ul	99
7) Acenaphthylene	14.67	152	3147	0.407	ng/ul	98
8) Acenaphthene	15.01	153	2239	0.401	ng/ul	93
9) Fluorene	15.97	166	2885	0.391	ng/ul	89
11) Pentachlorophenol	17.31	266	257	0.286	ng/ul	99
12) Phenanthrene	17.70	178	4592	0.417	ng/ul#	89
13) Anthracene	17.78	178	4362	0.395	ng/ul#	90
15) Fluoranthene	19.67	202	6390	0.380	ng/ul	85
17) Pyrene	20.02	202	2779	0.264	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	4207	0.265	ng/ul#	86
19) Chrysene	21.79	228	5970	0.409	ng/ul	97
21) Benzo(b)fluoranthene	23.54	252	5905	0.387	ng/ul	86
22) Benzo(k)fluoranthene	23.59	252	5234	0.359	ng/ul#	89
23) Benzo(a)pyrene	24.23	252	5035	0.380	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.13	276	5780	0.324	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.14	278	4702	0.314	ng/ul	93
26) Benzo(g,h,i)perylene	27.99	276	4837	0.311	ng/ul	97

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

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