

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042718\
 Data File : BE096202.D
 Acq On : 27 Apr 2018 11:02
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.468

Quant Time: Apr 27 23:51:58 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 04:45:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	773	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2812	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1533	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	3752	0.40	ng/ul	0.00
16) Chrysene-d12	21.74	240	5062	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3980	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2045	0.39	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	6067	0.35	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	11.23	128	3140	0.401	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	2111	0.402	ng/ul	99
7) Acenaphthylene	14.67	152	3256	0.412	ng/ul	98
8) Acenaphthene	15.01	153	2337	0.410	ng/ul	95
9) Fluorene	15.97	166	2612	0.346	ng/ul	91
11) Pentachlorophenol	17.31	266	283	0.313	ng/ul	92
12) Phenanthrene	17.70	178	4787	0.433	ng/ul#	89
13) Anthracene	17.78	178	4479	0.404	ng/ul#	91
15) Fluoranthene	19.67	202	6480	0.384	ng/ul	84
17) Pyrene	20.02	202	3291	0.324	ng/ul#	80
18) Benzo(a)anthracene	21.73	228	5723	0.373	ng/ul#	85
19) Chrysene	21.79	228	5731	0.407	ng/ul	97
21) Benzo(b)fluoranthene	23.54	252	5603	0.372	ng/ul	89
22) Benzo(k)fluoranthene	23.59	252	5189	0.360	ng/ul	91
23) Benzo(a)pyrene	24.23	252	5078	0.388	ng/ul#	84
24) Indeno(1,2,3-cd)pyrene	27.13	276	5603	0.318	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.15	278	4571	0.309	ng/ul	92
26) Benzo(g,h,i)perylene	28.00	276	4644	0.302	ng/ul	97

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

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