

Data Path : Z:\HPCHEM1\BNA E\DATA\BE042618\
 Data File : BE096200.D
 Acq On : 27 Apr 2018 3:45
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.467

Quant Time: Apr 27 05:05:05 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	675	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	1611	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.65	188	3761	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	4520	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	3788	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2091	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	5592	0.32	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.23	128	3110	0.397	ng/ul#	90
5) 2-Methylnaphthalene	12.81	142	2178	0.415	ng/ul	100
7) Acenaphthylene	14.67	152	3401	0.410	ng/ul	99
8) Acenaphthene	15.01	153	2429	0.406	ng/ul	93
9) Fluorene	15.97	166	2714	0.343	ng/ul	91
11) Pentachlorophenol	17.32	266	200	0.221	ng/ul	94
12) Phenanthrene	17.70	178	5302	0.479	ng/ul#	89
13) Anthracene	17.79	178	4477	0.403	ng/ul#	92
15) Fluoranthene	19.67	202	6970	0.412	ng/ul	85
17) Pyrene	20.02	202	3536	0.390	ng/ul#	80
18) Benzo(a)anthracene	21.73	228	5320	0.389	ng/ul#	85
19) Chrysene	21.79	228	5594	0.445	ng/ul	97
21) Benzo(b)fluoranthene	23.54	252	5307	0.370	ng/ul	87
22) Benzo(k)fluoranthene	23.60	252	4992	0.364	ng/ul#	89
23) Benzo(a)pyrene	24.24	252	4788	0.384	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.13	276	5457	0.325	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.15	278	4417	0.314	ng/ul	93
26) Benzo(g,h,i)perylene	28.00	276	4514	0.309	ng/ul	96

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

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