

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031418\
 Data File : BE095767.D
 Acq On : 14 Mar 2018 16:58
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.417

Quant Time: Mar 15 01:37:50 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 15 01:35:18 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1998	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	5807	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.01	164	3380	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	7489	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7725	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	7418	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3571	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	9327	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.30	128	5864	0.362	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	4194	0.372	ng/ul	99
7) Acenaphthylene	14.73	152	6192	0.363	ng/ul	97
8) Acenaphthene	15.07	153	4571	0.356	ng/ul	93
9) Fluorene	16.03	166	4949	0.355	ng/ul#	89
11) Pentachlorophenol	17.36	266	388	0.369	ng/ul	91
12) Phenanthrene	17.75	178	8078	0.360	ng/ul#	88
13) Anthracene	17.84	178	7896	0.377	ng/ul#	91
15) Fluoranthene	19.72	202	10240	0.377	ng/ul	84
17) Pyrene	20.07	202	11509	0.392	ng/ul#	77
18) Benzo(a)anthracene	21.78	228	9447	0.379	ng/ul#	86
19) Chrysene	21.84	228	9191	0.362	ng/ul	96
21) Benzo(b)fluoranthene	23.61	252	8759	0.341	ng/ul	85
22) Benzo(k)fluoranthene	23.66	252	8526	0.349	ng/ul#	87
23) Benzo(a)pyrene	24.30	252	8290	0.353	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.23	276	9266	0.356	ng/ul#	76
25) Dibenzo(a,h)anthracene	27.25	278	7489	0.360	ng/ul#	89
26) Benzo(g,h,i)perylene	28.10	276	7798	0.349	ng/ul#	93

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

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