

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE062417\
 Data File : BE093366.D
 Acq On : 24 Jun 2017 17:45
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.425

Quant Time: Jun 26 01:15:37 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jun 26 01:13:30 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	1545	0.40	ng/ul	0.00
2) Naphthalene-d8	10.51	136	7813	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.36	164	4533	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.10	188	11803	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	13302	0.40	ng/ul	0.00
20) Perylene-d12	23.66	264	11243	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.10	152	4742	0.41	ng/ul	0.00
14) Fluoranthene-d10	19.12	212	13195	0.41	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.56	128	7707	0.398	ng/ul#	92
5) 2-Methylnaphthalene	12.18	142	5515	0.416	ng/ul	98
7) Acenaphthylene	14.08	152	8346	0.429	ng/ul	97
8) Acenaphthene	14.42	153	6185	0.394	ng/ul	99
9) Fluorene	15.42	166	7123	0.387	ng/ul	96
11) Pentachlorophenol	16.76	266	821	0.469	ng/ul	98
12) Phenanthrene	17.14	178	12765	0.393	ng/ul#	91
13) Anthracene	17.22	178	12697	0.449	ng/ul#	91
15) Fluoranthene	19.15	202	15984	0.401	ng/ul	84
17) Pyrene	19.51	202	17284	0.441	ng/ul#	89
18) Benzo(a)anthracene	21.24	228	15357	0.433	ng/ul#	85
19) Chrysene	21.29	228	14327	0.389	ng/ul	98
21) Benzo(b)fluoranthene	22.92	252	14074	0.412	ng/ul	89
22) Benzo(k)fluoranthene	22.97	252	13664	0.377	ng/ul#	90
23) Benzo(a)pyrene	23.55	252	13297	0.412	ng/ul#	85
24) Indeno(1,2,3-cd)pyrene	26.18	276	15159	0.406	ng/ul#	93
25) Dibenzo(a,h)anthracene	26.20	278	12663	0.403	ng/ul#	86
26) Benzo(g,h,i)perylene	26.95	276	12511	0.391	ng/ul	96

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

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