

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE072417\
 Data File : BE093532.D
 Acq On : 24 Jul 2017 11:20
 Operator : SJ/JU
 Sample : SSTD0.832
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.832

Quant Time: Jul 24 11:53:21 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE072417.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 24 11:37:06 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.93	152	2685	0.40	ng/ul	0.00
2) Naphthalene-d8	10.72	136	13317	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.54	164	8521	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	23474	0.40	ng/ul	0.00
16) Chrysene-d12	21.42	240	29956	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	26458	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.30	152	17399	0.89	ng/ul	0.00
14) Fluoranthene-d10	19.28	212	60797	0.95	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	10.76	128	26229	0.794	ng/ul#	91
5) 2-Methylnaphthalene	12.37	142	18941	0.838	ng/ul	98
7) Acenaphthylene	14.26	152	28547	0.781	ng/ul	96
8) Acenaphthene	14.60	153	22255	0.754	ng/ul	97
9) Fluorene	15.58	166	28418	0.822	ng/ul#	94
11) Pentachlorophenol	16.93	266	2947	0.846	ng/ul	96
12) Phenanthrene	17.31	178	50847	0.787	ng/ul#	93
13) Anthracene	17.39	178	50427	0.897	ng/ul#	91
15) Fluoranthene	19.31	202	64612	0.816	ng/ul	83
17) Pyrene	19.67	202	67098	0.760	ng/ul#	83
18) Benzo(a)anthracene	21.40	228	67960	0.851	ng/ul#	86
19) Chrysene	21.46	228	65928	0.795	ng/ul	98
21) Benzo(b)fluoranthene	23.15	252	65050	0.810	ng/ul	84
22) Benzo(k)fluoranthene	23.20	252	64079	0.751	ng/ul#	87
23) Benzo(a)pyrene	23.81	252	61371	0.808	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	26.57	276	67888	0.772	ng/ul#	83
25) Dibenzo(a,h)anthracene	26.59	278	57428	0.776	ng/ul#	88
26) Benzo(g,h,i)perylene	27.38	276	57262	0.760	ng/ul	95

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

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