

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095880.D
 Acq On : 28 Mar 2018 17:09
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.424

Quant Time: Mar 29 04:33:32 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Mar 28 16:31:37 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.41	152	1589	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	4158	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.99	164	2525	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	6710	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	7385	0.40	ng/ul	0.00
20) Perylene-d12	24.42	264	7088	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.79	152	2963	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	10003	0.47	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	4245	0.366	ng/ul#	89
5) 2-Methylnaphthalene	12.87	142	2976	0.369	ng/ul	99
7) Acenaphthylene	14.72	152	4847	0.380	ng/ul	98
8) Acenaphthene	15.06	153	3513	0.366	ng/ul	95
9) Fluorene	16.02	166	4094	0.393	ng/ul#	91
11) Pentachlorophenol	17.36	266	322	0.341	ng/ul	91
12) Phenanthrene	17.74	178	7285	0.363	ng/ul#	88
13) Anthracene	17.83	178	7092	0.377	ng/ul#	90
15) Fluoranthene	19.71	202	9729	0.400	ng/ul	84
17) Pyrene	20.06	202	10840	0.387	ng/ul#	79
18) Benzo(a)anthracene	21.77	228	9193	0.386	ng/ul#	84
19) Chrysene	21.83	228	8662	0.357	ng/ul	97
21) Benzo(b)fluoranthene	23.59	252	8377	0.341	ng/ul	87
22) Benzo(k)fluoranthene	23.65	252	8423	0.361	ng/ul#	89
23) Benzo(a)pyrene	24.29	252	8119	0.362	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	27.21	276	9150	0.368	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.23	278	7422	0.373	ng/ul	92
26) Benzo(g,h,i)perylene	28.09	276	7615	0.357	ng/ul#	94

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

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