

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE042518\
 Data File : BE096150.D
 Acq On : 25 Apr 2018 16:25
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.462

Quant Time: Apr 26 04:20:11 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE042418.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Apr 26 04:18:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.36	152	725	0.40	ng/ul	0.00
2) Naphthalene-d8	11.18	136	2724	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.95	164	1650	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.66	188	4728	0.40	ng/ul	0.00
16) Chrysene-d12	21.75	240	6284	0.40	ng/ul	0.00
20) Perylene-d12	24.35	264	5263	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.74	152	2024	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.64	212	8032	0.37	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	11.23	128	3067	0.404	ng/ul#	91
5) 2-Methylnaphthalene	12.81	142	2109	0.415	ng/ul	98
7) Acenaphthylene	14.67	152	3382	0.398	ng/ul	98
8) Acenaphthene	15.01	153	2936	0.479	ng/ul	92
9) Fluorene	15.97	166	4258	0.525	ng/ul	87
11) Pentachlorophenol	17.31	266	378	0.332	ng/ul	97
12) Phenanthrene	17.70	178	6688	0.480	ng/ul#	89
13) Anthracene	17.79	178	5696	0.408	ng/ul#	92
15) Fluoranthene	19.67	202	8348	0.392	ng/ul	85
17) Pyrene	20.02	202	4541	0.360	ng/ul#	79
18) Benzo(a)anthracene	21.73	228	6381	0.335	ng/ul#	86
19) Chrysene	21.79	228	7022	0.402	ng/ul	97
21) Benzo(b)fluoranthene	23.55	252	7208	0.362	ng/ul	87
22) Benzo(k)fluoranthene	23.60	252	6638	0.348	ng/ul#	89
23) Benzo(a)pyrene	24.24	252	6590	0.381	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	27.13	276	7525	0.323	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.14	278	6189	0.317	ng/ul	91
26) Benzo(g,h,i)perylene	28.00	276	6356	0.313	ng/ul	97

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

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