

Data Path : Z:\HPCHEM1\BNA E\DATA\BE031518\
 Data File : BE095806.D
 Acq On : 15 Mar 2018 16:40
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.421

Quant Time: Mar 16 01:14:21 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Mar 16 01:12:14 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.45	152	2154	0.40	ng/ul	0.00
2) Naphthalene-d8	11.28	136	6122	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.04	164	3243	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.75	188	6852	0.40	ng/ul	0.00
16) Chrysene-d12	21.84	240	7027	0.40	ng/ul	0.00
20) Perylene-d12	24.50	264	7000	0.40	ng/ul	0.01

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.83	152	3677	0.35	ng/ul	0.00
14) Fluoranthene-d10	19.72	212	8126	0.37	ng/ul	0.00

Target Compounds				Ovalue		
3) Naphthalene	11.32	128	6183	0.362	ng/ul#	90
5) 2-Methylnaphthalene	12.91	142	4202	0.354	ng/ul	99
7) Acenaphthylene	14.76	152	6318	0.386	ng/ul	97
8) Acenaphthene	15.10	153	4410	0.358	ng/ul	95
9) Fluorene	16.06	166	4725	0.353	ng/ul	94
11) Pentachlorophenol	17.41	266	354	0.368	ng/ul#	85
12) Phenanthrene	17.78	178	7258	0.354	ng/ul	92
13) Anthracene	17.87	178	7273	0.379	ng/ul	93
15) Fluoranthene	19.75	202	8957	0.360	ng/ul	83
17) Pyrene	20.10	202	9046	0.339	ng/ul#	78
18) Benzo(a)anthracene	21.81	228	8660	0.382	ng/ul#	78
19) Chrysene	21.87	228	8061	0.349	ng/ul#	74
21) Benzo(b)fluoranthene	23.67	252	8570	0.353	ng/ul#	63
22) Benzo(k)fluoranthene	23.72	252	7726	0.335	ng/ul#	59
23) Benzo(a)pyrene	24.38	252	7864	0.355	ng/ul#	60
24) Indeno(1,2,3-cd)pyrene	27.36	276	8328	0.339	ng/ul#	72
25) Dibenzo(a,h)anthracene	27.37	278	7178	0.366	ng/ul#	55
26) Benzo(g,h,i)perylene	28.26	276	6269	0.297	ng/ul#	56

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

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