

Data Path : Z:\HPCHEM1\BNA E\DATA\BE032818\
 Data File : BE095904.D
 Acq On : 29 Mar 2018 7:27
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.426

Quant Time: Mar 29 08:07:33 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE031318.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Mar 29 07:00:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.42	152	1752	0.40	ng/ul	0.00
2) Naphthalene-d8	11.24	136	5299	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.00	164	3133	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.70	188	7147	0.40	ng/ul	0.00
16) Chrysene-d12	21.79	240	6988	0.40	ng/ul	0.00
20) Perylene-d12	24.43	264	6919	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.80	152	3834	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.68	212	9817	0.43	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.29	128	5390	0.364	ng/ul#	88
5) 2-Methylnaphthalene	12.87	142	3861	0.375	ng/ul	99
7) Acenaphthylene	14.73	152	6234	0.394	ng/ul	96
8) Acenaphthene	15.06	153	4370	0.367	ng/ul	95
9) Fluorene	16.03	166	4848	0.375	ng/ul#	91
11) Pentachlorophenol	17.37	266	445	0.443	ng/ul	99
12) Phenanthrene	17.75	178	7586	0.354	ng/ul#	89
13) Anthracene	17.84	178	7757	0.388	ng/ul#	91
15) Fluoranthene	19.72	202	9393	0.362	ng/ul	84
17) Pyrene	20.07	202	10693	0.403	ng/ul#	77
18) Benzo(a)anthracene	21.78	228	8878	0.394	ng/ul#	87
19) Chrysene	21.84	228	8337	0.363	ng/ul	91
21) Benzo(b)fluoranthene	23.61	252	8186	0.342	ng/ul	92
22) Benzo(k)fluoranthene	23.66	252	8095	0.355	ng/ul#	94
23) Benzo(a)pyrene	24.31	252	7941	0.363	ng/ul	90
24) Indeno(1,2,3-cd)pyrene	27.24	276	9207	0.379	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.25	278	7405	0.382	ng/ul#	91
26) Benzo(g,h,i)perylene	28.12	276	7450	0.357	ng/ul#	92

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

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