

Data Path : Z:\HPCHEM1\BNA_E\DATA\BE040918\
 Data File : BE095961.D
 Acq On : 9 Apr 2018 9:36
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.436

Quant Time: Apr 10 03:43:15 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE040618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Apr 07 00:03:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.39	152	1758	0.40	ng/ul	0.00
2) Naphthalene-d8	11.21	136	4443	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.98	164	2535	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.68	188	6197	0.40	ng/ul	0.00
16) Chrysene-d12	21.77	240	7213	0.40	ng/ul	0.00
20) Perylene-d12	24.40	264	6854	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.77	152	3295	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.66	212	10132	0.40	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.26	128	4922	0.397	ng/ul#	89
5) 2-Methylnaphthalene	12.84	142	3367	0.400	ng/ul	99
7) Acenaphthylene	14.70	152	5142	0.391	ng/ul	98
8) Acenaphthene	15.04	153	3702	0.393	ng/ul	96
9) Fluorene	16.00	166	4203	0.395	ng/ul	88
11) Pentachlorophenol	17.34	266	382	0.291	ng/ul	98
12) Phenanthrene	17.72	178	7250	0.397	ng/ul#	89
13) Anthracene	17.81	178	7197	0.396	ng/ul#	92
15) Fluoranthene	19.69	202	9953	0.400	ng/ul	84
17) Pyrene	20.05	202	10885	0.444	ng/ul#	81
18) Benzo(a)anthracene	21.76	228	12776	0.438	ng/ul#	86
19) Chrysene	21.81	228	9323	0.445	ng/ul	97
21) Benzo(b)fluoranthene	23.58	252	9062	0.409	ng/ul	85
22) Benzo(k)fluoranthene	23.63	252	8675	0.393	ng/ul#	89
23) Benzo(a)pyrene	24.28	252	8504	0.397	ng/ul#	81
24) Indeno(1,2,3-cd)pyrene	27.19	276	9397	0.395	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.21	278	7560	0.395	ng/ul	91
26) Benzo(g,h,i)perylene	28.06	276	7982	0.402	ng/ul	95

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

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