

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095448.D
 Acq On : 3 Feb 2018 21:34
 Operator : SJ/JU
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.439

Quant Time: Feb 05 03:35:26 2018
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Feb 05 03:32:57 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2554	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7860	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4533	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10217	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10499	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9917	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5516	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	13622	0.35	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.31	128	9157	0.416	ng/ul#	89
5) 2-Methylnaphthalene	12.88	142	6413	0.436	ng/ul	99
7) Acenaphthylene	14.74	152	9485	0.435	ng/ul	97
8) Acenaphthene	15.08	153	6905	0.408	ng/ul	97
9) Fluorene	16.03	166	7936	0.455	ng/ul#	85
11) Pentachlorophenol	17.36	266	641	0.324	ng/ul	96
12) Phenanthrene	17.75	178	13215	0.385	ng/ul#	92
13) Anthracene	17.85	178	12080	0.384	ng/ul#	93
15) Fluoranthene	19.72	202	15535	0.349	ng/ul	85
17) Pyrene	20.07	202	16585	0.471	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	14405	0.439	ng/ul#	84
19) Chrysene	21.84	228	13954	0.409	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	13962	0.399	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	12679	0.396	ng/ul#	85
23) Benzo(a)pyrene	24.31	252	12842	0.411	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	27.24	276	13008	0.380	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.26	278	10343	0.369	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	11018	0.380	ng/ul	94

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

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