

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020318\
 Data File : BE095433.D
 Acq On : 3 Feb 2018 12:37
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.438

Quant Time: Feb 05 03:30:28 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 01:47:56 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2647	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	8115	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4642	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	10605	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	11009	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	10312	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5502	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	14007	0.35	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	11.31	128	9176	0.404	ng/ul#	90
5) 2-Methylnaphthalene	12.88	142	6336	0.418	ng/ul	100
7) Acenaphthylene	14.74	152	9444	0.423	ng/ul	97
8) Acenaphthene	15.08	153	7062	0.408	ng/ul	97
9) Fluorene	16.03	166	8190	0.458	ng/ul#	87
11) Pentachlorophenol	17.36	266	715	0.348	ng/ul	98
12) Phenanthrene	17.75	178	13315	0.373	ng/ul#	91
13) Anthracene	17.85	178	12238	0.375	ng/ul#	92
15) Fluoranthene	19.72	202	15890	0.344	ng/ul	83
17) Pyrene	20.07	202	17386	0.471	ng/ul#	80
18) Benzo(a)anthracene	21.78	228	14666	0.426	ng/ul#	83
19) Chrysene	21.84	228	14380	0.402	ng/ul	97
21) Benzo(b)fluoranthene	23.62	252	14321	0.394	ng/ul	83
22) Benzo(k)fluoranthene	23.67	252	13351	0.401	ng/ul#	85
23) Benzo(a)pyrene	24.31	252	13291	0.409	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	14155	0.397	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.26	278	11235	0.386	ng/ul#	87
26) Benzo(g,h,i)perylene	28.12	276	11998	0.398	ng/ul#	94

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

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