

Data Path : \\74.0.250.170\SV0ASRV\HPCHEM1\BNA_E\DATA\BE020518\
 Data File : BE095484.D
 Acq On : 5 Feb 2018 10:58
 Operator : SJ/JU
 Sample : SSTDCCC0.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.442

Quant Time: Feb 06 01:22:54 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 04:46:10 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.43	152	1530	0.40	ng/ul	0.00
2) Naphthalene-d8	11.25	136	3851	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	2268	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	5962	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	6442	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	6033	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	2624	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.69	212	8006	0.35	ng/ul	0.00
Target Compounds					Qvalue	
3) Naphthalene	11.30	128	4427	0.411	ng/ul#	90
5) 2-Methylnaphthalene	12.88	142	3028	0.421	ng/ul	100
7) Acenaphthylene	14.74	152	4553	0.417	ng/ul	98
8) Acenaphthene	15.08	153	3406	0.402	ng/ul	97
9) Fluorene	16.03	166	4169	0.478	ng/ul#	89
11) Pentachlorophenol	17.36	266	373	0.323	ng/ul	99
12) Phenanthrene	17.75	178	7087	0.353	ng/ul#	91
13) Anthracene	17.85	178	6431	0.350	ng/ul	94
15) Fluoranthene	19.72	202	9021	0.347	ng/ul	83
17) Pyrene	20.07	202	9817	0.454	ng/ul#	81
18) Benzo(a)anthracene	21.78	228	8619	0.428	ng/ul#	84
19) Chrysene	21.84	228	8377	0.401	ng/ul	98
21) Benzo(b)fluoranthene	23.62	252	7956	0.374	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	7631	0.392	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	7314	0.385	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	7127	0.342	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.26	278	5487	0.322	ng/ul	92
26) Benzo(g,h,i)perylene	28.12	276	6474	0.367	ng/ul	96

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

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