

Data Path : Z:\HPCHEM1\BNA_E\Data\BE092017\
 Data File : BE094062.D
 Acq On : 21 Sep 2017 9:49
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.411

Quant Time: Sep 21 11:05:23 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE092017.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 21 07:18:34 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.91	152	1762	0.40	ng/ul	0.00
2) Naphthalene-d8	10.70	136	7709	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.53	164	3990	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.26	188	9572	0.40	ng/ul	-0.02
16) Chrysene-d12	21.42	240	12226	0.40	ng/ul	0.00
20) Perylene-d12	23.92	264	12518	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.28	152	4626	0.39	ng/ul	-0.02
14) Fluoranthene-d10	19.28	212	12294	0.38	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.75	128	7466	0.395	ng/ul#	93
5) 2-Methylnaphthalene	12.35	142	4829	0.394	ng/ul	100
7) Acenaphthylene	14.25	152	6675	0.429	ng/ul	98
8) Acenaphthene	14.59	153	5113	0.399	ng/ul	98
9) Fluorene	15.58	166	5647	0.377	ng/ul	95
11) Pentachlorophenol	16.94	266	426	0.379	ng/ul	97
12) Phenanthrene	17.31	178	9730	0.389	ng/ul#	92
13) Anthracene	17.39	178	9442	0.388	ng/ul#	93
15) Fluoranthene	19.31	202	12494	0.380	ng/ul	84
17) Pyrene	19.67	202	13862	0.408	ng/ul#	82
18) Benzo(a)anthracene	21.40	228	13853	0.427	ng/ul#	86
19) Chrysene	21.45	228	15069	0.402	ng/ul	97
21) Benzo(b)fluoranthene	23.15	252	13507	0.397	ng/ul	88
22) Benzo(k)fluoranthene	23.20	252	15036	0.383	ng/ul#	89
23) Benzo(a)pyrene	23.81	252	14192	0.402	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	26.59	276	15024	0.416	ng/ul#	84
25) Dibenzo(a,h)anthracene	26.60	278	12581	0.415	ng/ul#	92
26) Benzo(g,h,i)perylene	27.39	276	12683	0.407	ng/ul	99

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

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