

Data Path : Z:\HPCHEM1\BNA_E\Data\BE063017\
 Data File : BE093375.D
 Acq On : 30 Jun 2017 10:23
 Operator : SJ/JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.428

Quant Time: Jun 30 11:12:33 2017
 Quant Method : Z:\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE060717.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jun 30 10:50:58 2017
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.95	152	3961	0.40	ng/ul	0.00
2) Naphthalene-d8	10.74	136	19661	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.56	164	12079	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.27	188	33372	0.40	ng/ul	-0.02
16) Chrysene-d12	21.44	240	47456	0.40	ng/ul	0.00
20) Perylene-d12	23.95	264	45616	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.32	152	12320	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.30	212	42886	0.47	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.79	128	20084	0.412	ng/ul#	92
5) 2-Methylnaphthalene	12.39	142	14340	0.430	ng/ul	99
7) Acenaphthylene	14.27	152	22071	0.426	ng/ul#	96
8) Acenaphthene	14.62	153	16467	0.394	ng/ul	98
9) Fluorene	15.60	166	20189	0.412	ng/ul#	94
11) Pentachlorophenol	16.95	266	1120	0.226	ng/ul	96
12) Phenanthrene	17.33	178	36797	0.401	ng/ul#	92
13) Anthracene	17.41	178	35514	0.444	ng/ul#	92
15) Fluoranthene	19.33	202	48463	0.430	ng/ul	84
17) Pyrene	19.69	202	52456	0.375	ng/ul#	83
18) Benzo(a)anthracene	21.42	228	53741	0.425	ng/ul#	85
19) Chrysene	21.48	228	52720	0.401	ng/ul	97
21) Benzo(b)fluoranthene	23.19	252	54682	0.395	ng/ul	84
22) Benzo(k)fluoranthene	23.23	252	57121	0.389	ng/ul#	89
23) Benzo(a)pyrene	23.84	252	51756	0.395	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	26.62	276	60129	0.397	ng/ul#	81
25) Dibenzo(a,h)anthracene	26.64	278	50892	0.399	ng/ul#	88
26) Benzo(g,h,i)perylene	27.43	276	51163	0.394	ng/ul	94

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

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