

Data Path : Z:\HPCHEM1\BNA E\DATA\BE012818\
 Data File : BE095205.D
 Acq On : 28 Jan 2018 14:04
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.417

Quant Time: Jan 29 06:56:43 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 Jan 29 01:25:04 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2533	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7588	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4494	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9936	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	11356	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	10550	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.82	152	5163	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	14749	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8438	0.397	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5923	0.417	ng/ul	100
7) Acenaphthylene	14.74	152	9145	0.423	ng/ul	96
8) Acenaphthene	15.08	153	6824	0.407	ng/ul	97
9) Fluorene	16.02	166	7269	0.420	ng/ul#	75
11) Pentachlorophenol	17.36	266	907	0.471	ng/ul	95
12) Phenanthrene	17.75	178	14164	0.424	ng/ul#	87
13) Anthracene	17.85	178	13699	0.448	ng/ul#	91
15) Fluoranthene	19.72	202	16627	0.384	ng/ul	85
17) Pyrene	20.08	202	18571	0.487	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	15352	0.432	ng/ul#	85
19) Chrysene	21.84	228	14807	0.402	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	14001	0.376	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	13891	0.408	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	13431	0.404	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.24	276	15050	0.413	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.27	278	12322	0.413	ng/ul#	90
26) Benzo(g,h,i)perylene	28.12	276	12609	0.409	ng/ul	96

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

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