

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095301.D
 Acq On : 31 Jan 2018 5:25
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.425

Quant Time: Jan 31 07:51:46 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Jan 31 03:40:45 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2495	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7688	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4471	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9327	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10102	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9439	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5246	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13104	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8555	0.397	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	6027	0.419	ng/ul	99
7) Acenaphthylene	14.75	152	9023	0.419	ng/ul	96
8) Acenaphthene	15.08	153	6683	0.400	ng/ul	94
9) Fluorene	16.02	166	7094	0.412	ng/ul#	75
11) Pentachlorophenol	17.36	266	762	0.422	ng/ul	99
12) Phenanthrene	17.75	178	12680	0.404	ng/ul#	86
13) Anthracene	17.84	178	12395	0.432	ng/ul#	92
15) Fluoranthene	19.72	202	14622	0.360	ng/ul	84
17) Pyrene	20.08	202	16609	0.490	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	13600	0.431	ng/ul#	86
19) Chrysene	21.84	228	13073	0.399	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12812	0.385	ng/ul	84
22) Benzo(k)fluoranthene	23.67	252	12452	0.409	ng/ul#	86
23) Benzo(a)pyrene	24.31	252	12282	0.413	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	13525	0.415	ng/ul#	80
25) Dibenzo(a,h)anthracene	27.27	278	10907	0.409	ng/ul#	88
26) Benzo(g,h,i)perylene	28.12	276	11374	0.412	ng/ul	95

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

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