

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095291.D
 Acq On : 30 Jan 2018 18:26
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.424

Quant Time: Jan 31 01:27:03 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 01:22:09 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2372	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7272	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4273	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	8822	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9902	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9389	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4978	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13005	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8086	0.397	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5764	0.424	ng/ul	100
7) Acenaphthylene	14.74	152	8786	0.427	ng/ul	96
8) Acenaphthene	15.08	153	6439	0.404	ng/ul	94
9) Fluorene	16.03	166	6826	0.415	ng/ul#	76
11) Pentachlorophenol	17.36	266	845	0.495	ng/ul	95
12) Phenanthrene	17.75	178	12468	0.420	ng/ul#	88
13) Anthracene	17.85	178	12090	0.445	ng/ul#	92
15) Fluoranthene	19.73	202	14427	0.376	ng/ul	83
17) Pyrene	20.08	202	16830	0.507	ng/ul#	79
18) Benzo(a)anthracene	21.79	228	13495	0.436	ng/ul#	86
19) Chrysene	21.84	228	12786	0.398	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12450	0.376	ng/ul	87
22) Benzo(k)fluoranthene	23.67	252	12280	0.405	ng/ul#	88
23) Benzo(a)pyrene	24.32	252	11935	0.404	ng/ul#	80
24) Indeno(1,2,3-cd)pyrene	27.24	276	13474	0.415	ng/ul#	78
25) Dibenzo(a,h)anthracene	27.27	278	10917	0.411	ng/ul#	90
26) Benzo(g,h,i)perylene	28.12	276	11264	0.410	ng/ul	94

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

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