

Data Path : Z:\HPCHEM1\BNA E\DATA\BE013018\
 Data File : BE095309.D
 Acq On : 31 Jan 2018 10:12
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.426

Quant Time: Jan 31 11:47:13 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 08:16:13 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2247	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	6978	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4180	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9445	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	10021	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9628	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	4792	0.42	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	13356	0.37	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	7799	0.399	ng/ul#	89
5) 2-Methylnaphthalene	12.89	142	5559	0.426	ng/ul	99
7) Acenaphthylene	14.75	152	8679	0.431	ng/ul	96
8) Acenaphthene	15.08	153	6355	0.407	ng/ul	95
9) Fluorene	16.03	166	7123	0.443	ng/ul#	81
11) Pentachlorophenol	17.36	266	925	0.506	ng/ul	97
12) Phenanthrene	17.75	178	12816	0.403	ng/ul#	90
13) Anthracene	17.85	178	12372	0.426	ng/ul#	93
15) Fluoranthene	19.72	202	14854	0.361	ng/ul	84
17) Pyrene	20.08	202	16948	0.504	ng/ul#	77
18) Benzo(a)anthracene	21.79	228	13663	0.436	ng/ul#	86
19) Chrysene	21.84	228	13015	0.400	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12825	0.378	ng/ul	87
22) Benzo(k)fluoranthene	23.67	252	12292	0.396	ng/ul#	88
23) Benzo(a)pyrene	24.31	252	12174	0.401	ng/ul#	82
24) Indeno(1,2,3-cd)pyrene	27.24	276	13627	0.410	ng/ul#	79
25) Dibenzo(a,h)anthracene	27.26	278	11148	0.410	ng/ul#	90
26) Benzo(g,h,i)perylene	28.11	276	11396	0.405	ng/ul	93

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

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