

Data Path : Z:\HPCHEM1\BNA E\DATA\BE020118\
 Data File : BE095357.D
 Acq On : 1 Feb 2018 15:06
 Operator : SJ/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.431

Quant Time: Feb 02 04:04:51 2018
 Quant Method : Z:\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE012618.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Feb 01 06:27:03 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	8.44	152	2353	0.40	ng/ul	0.00
2) Naphthalene-d8	11.26	136	7405	0.40	ng/ul	0.00
6) Acenaphthene-d10	15.02	164	4310	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.71	188	9066	0.40	ng/ul	0.00
16) Chrysene-d12	21.80	240	9558	0.40	ng/ul	0.00
20) Perylene-d12	24.44	264	9167	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.81	152	5147	0.43	ng/ul	0.00
14) Fluoranthene-d10	19.70	212	12380	0.36	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	11.31	128	8321	0.401	ng/ul#	88
5) 2-Methylnaphthalene	12.89	142	5937	0.429	ng/ul	100
7) Acenaphthylene	14.75	152	8945	0.431	ng/ul	96
8) Acenaphthene	15.08	153	6496	0.404	ng/ul	95
9) Fluorene	16.03	166	6946	0.419	ng/ul#	81
11) Pentachlorophenol	17.36	266	628	0.358	ng/ul	96
12) Phenanthrene	17.75	178	11993	0.393	ng/ul#	90
13) Anthracene	17.85	178	11465	0.411	ng/ul#	93
15) Fluoranthene	19.72	202	13848	0.351	ng/ul	84
17) Pyrene	20.08	202	15283	0.477	ng/ul#	78
18) Benzo(a)anthracene	21.79	228	12980	0.434	ng/ul#	86
19) Chrysene	21.84	228	12302	0.396	ng/ul	96
21) Benzo(b)fluoranthene	23.62	252	12204	0.377	ng/ul	85
22) Benzo(k)fluoranthene	23.67	252	11764	0.398	ng/ul#	87
23) Benzo(a)pyrene	24.32	252	11538	0.400	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	27.25	276	12835	0.405	ng/ul#	77
25) Dibenzo(a,h)anthracene	27.27	278	10333	0.399	ng/ul#	90
26) Benzo(g,h,i)perylene	28.13	276	10778	0.402	ng/ul#	93

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

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