

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE081718\
 Data File : BE097294.D
 Acq On : 17 Aug 2018 12:54
 Operator : SJ/JU
 Sample : SSTD1.605
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD1.605

Quant Time: Aug 17 13:32:57 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE081718.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 17 12:19:40 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	913	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	4679	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2525	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	5601	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	5526	0.40	ng/ul	0.00
20) Perylene-d12	23.22	264	4276	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	12424	1.65	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	27031	1.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.18	128	18543	1.637	ng/ul#	91
5) 2-Methylnaphthalene	11.80	142	12619	1.686	ng/ul	99
7) Acenaphthylene	13.73	152	18691	1.712	ng/ul	96
8) Acenaphthene	14.08	153	13556	1.717	ng/ul#	91
9) Fluorene	15.08	166	15238	1.679	ng/ul#	93
11) Pentachlorophenol	16.44	266	679	0.846	ng/ul	96
12) Phenanthrene	16.81	178	24589	1.722	ng/ul#	89
13) Anthracene	16.90	178	22621	1.698	ng/ul#	92
15) Fluoranthene	18.84	202	28118	1.490	ng/ul	84
17) Pyrene	19.20	202	25381	1.714	ng/ul#	80
18) Benzo(a)anthracene	20.96	228	24281	1.538	ng/ul#	84
19) Chrysene	21.01	228	26289	1.616	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	22433	1.925	ng/ul	82
22) Benzo(k)fluoranthene	22.58	252	24151	1.922	ng/ul#	83
23) Benzo(a)pyrene	23.12	252	19781	1.696	ng/ul#	76
24) Indeno(1,2,3-cd)pyrene	25.52	276	23577	1.615	ng/ul#	76
25) Dibenzo(a,h)anthracene	25.53	278	19687	1.648	ng/ul#	85
26) Benzo(g,h,i)perylene	26.21	276	18767	1.488	ng/ul#	91

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

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