

Data Path : Z:\SVOASRV\HPCHEM1\BNA E\DATA\BE080218\
 Data File : BE097170.D
 Acq On : 3 Aug 2018 3:01
 Operator : SJ/JU
 Sample : SSTDC000.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD0.484

Quant Time: Aug 03 04:32:29 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA E\METHODS\SOM-EPA-SIM-BE080118.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 02 23:40:30 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.39	152	1342	0.40	ng/ul	0.00
2) Naphthalene-d8	10.13	136	6610	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.02	164	4081	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.77	188	10965	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	11864	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	12027	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.73	152	4364	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	13799	0.37	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.18	128	6350	0.397	ng/ul#	93
5) 2-Methylnaphthalene	11.80	142	4288	0.405	ng/ul	100
7) Acenaphthylene	13.73	152	7007	0.397	ng/ul	97
8) Acenaphthene	14.08	153	5092	0.399	ng/ul	92
9) Fluorene	15.08	166	5882	0.401	ng/ul#	93
11) Pentachlorophenol	16.43	266	646	0.411	ng/ul	97
12) Phenanthrene	16.81	178	11008	0.394	ng/ul#	89
13) Anthracene	16.90	178	10522	0.404	ng/ul#	92
15) Fluoranthene	18.84	202	13490	0.365	ng/ul	85
17) Pyrene	19.20	202	13968	0.439	ng/ul#	79
18) Benzo(a)anthracene	20.96	228	13621	0.402	ng/ul#	85
19) Chrysene	21.01	228	13748	0.394	ng/ul	96
21) Benzo(b)fluoranthene	22.53	252	12730	0.388	ng/ul	83
22) Benzo(k)fluoranthene	22.58	252	13937	0.394	ng/ul#	90
23) Benzo(a)pyrene	23.11	252	12814	0.391	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.50	276	15895	0.387	ng/ul#	78
25) Dibenzo(a,h)anthracene	25.52	278	13268	0.395	ng/ul#	89
26) Benzo(g,h,i)perylene	26.19	276	13494	0.380	ng/ul#	92

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

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