

Data Path : Z:\svoasrv\HPCHEM1\BNA_E\Data\BE082918\
 Data File : BE097386.D
 Acq On : 29 Aug 2018 12:29
 Operator : SJ/JU
 Sample : SSTD3.218
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_E
 ClientSampleId :
 SSTD3.218

Quant Time: Aug 29 13:20:46 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_E\METHODS\SOM-EPA-SIM-BE082918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Aug 29 13:07:59 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.38	152	939	0.40	ng/ul	0.00
2) Naphthalene-d8	10.12	136	4855	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.01	164	2947	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	7807	0.40	ng/ul	0.00
16) Chrysene-d12	20.97	240	10115	0.40	ng/ul	0.00
20) Perylene-d12	23.20	264	8535	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.72	152	22862	2.94	ng/ul	0.00
14) Fluoranthene-d10	18.80	212	74262	3.55	ng/ul	0.00

Target Compounds					Qvalue
3) Naphthalene	10.17	128	33737	2.848	ng/ul# 90
5) 2-Methylnaphthalene	11.79	142	23454	3.070	ng/ul 97
7) Acenaphthylene	13.72	152	33421	2.511	ng/ul 96
8) Acenaphthene	14.07	153	26231	2.739	ng/ul 95
9) Fluorene	15.07	166	31218	2.996	ng/ul# 93
11) Pentachlorophenol	16.42	266	3132	6.181	ng/ul 97
12) Phenanthrene	16.80	178	56651	2.808	ng/ul# 88
13) Anthracene	16.89	178	54458	3.222	ng/ul# 90
15) Fluoranthene	18.83	202	74905	3.495	ng/ul 84
17) Pyrene	19.19	202	72516	2.373	ng/ul# 80
18) Benzo(a)anthracene	20.95	228	78039	3.112	ng/ul# 85
19) Chrysene	21.00	228	80163	2.678	ng/ul 97
21) Benzo(b)fluoranthene	22.53	252	72275	3.127	ng/ul 81
22) Benzo(k)fluoranthene	22.57	252	74913	2.611	ng/ul# 83
23) Benzo(a)pyrene	23.10	252	67048	3.141	ng/ul# 75
24) Indeno(1,2,3-cd)pyrene	25.48	276	82653	3.113	ng/ul# 75
25) Dibenzo(a,h)anthracene	25.51	278	70102	3.264	ng/ul# 84
26) Benzo(g,h,i)perylene	26.18	276	70031	3.293	ng/ul# 89

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

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