

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN120518\
 Data File : BN003728.D
 Acq On : 06 Dec 2018 11:36
 Operator : JU/SJ
 Sample : SSTDCCC0.4EC
 Misc :
 ALS Vial : 22 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.488

Quant Time: Dec 06 12:17:37 2018
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN111918.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Dec 06 02:51:43 2018
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.83	152	7183	0.40	ng/ul	0.00
2) Naphthalene-d8	10.63	136	28214	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.47	164	14222	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.22	188	30986	0.40	ng/ul	0.00
16) Chrysene-d12	21.40	240	27633	0.40	ng/ul	0.00
20) Perylene-d12	23.71	264	23862	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.22	152	16637	0.38	ng/ul	0.00
14) Fluoranthene-d10	19.24	212	34237	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.68	128	33013	0.417	ng/ul	100
5) 2-Methylnaphthalene	12.29	142	22296	0.403	ng/ul	100
7) Acenaphthylene	14.19	152	26190	0.409	ng/ul	100
8) Acenaphthene	14.53	153	23135	0.412	ng/ul	100
9) Fluorene	15.52	166	25686	0.393	ng/ul	99
11) Pentachlorophenol	16.86	266	2551	0.457	ng/ul	99
12) Phenanthrene	17.26	178	42132	0.414	ng/ul	99
13) Anthracene	17.35	178	34664	0.415	ng/ul	100
15) Fluoranthene	19.27	202	47322	0.394	ng/ul	97
17) Pyrene	19.64	202	46685	0.471	ng/ul	99
18) Benzo(a)anthracene	21.39	228	37670	0.432	ng/ul	99
19) Chrysene	21.44	228	40159	0.402	ng/ul	100
21) Benzo(b)fluoranthene	23.01	252	39930	0.405	ng/ul	99
22) Benzo(k)fluoranthene	23.06	252	39403	0.410	ng/ul	99
23) Benzo(a)pyrene	23.62	252	35825	0.419	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	26.06	276	39020	0.398	ng/ul#	98
25) Dibenzo(a,h)anthracene	26.08	278	30785	0.390	ng/ul	99
26) Benzo(g,h,i)perylene	26.79	276	32941	0.387	ng/ul	99

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

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