

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN071519\
 Data File : BN006839.D
 Acq On : 15 Jul 2019 09:07
 Operator : HP/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.427

Quant Time: Jul 15 10:48:47 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN062719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jul 12 06:39:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.59	152	2214	0.40	ng/ul	0.00
2) Naphthalene-d8	10.36	136	8218	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	4244	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	9107	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	7789	0.40	ng/ul	0.00
20) Perylene-d12	23.46	264	9086	0.40	ng/ul	0.00

System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	4590	0.36	ng/ul	0.00
14) Fluoranthene-d10	19.05	212	10114	0.38	ng/ul	0.00

Target Compounds				Qvalue		
3) Naphthalene	10.41	128	9192	0.390	ng/ul	97
5) 2-Methylnaphthalene	12.04	142	5616	0.352	ng/ul	98
7) Acenaphthylene	13.96	152	7938	0.362	ng/ul	98
8) Acenaphthene	14.30	153	6413	0.380	ng/ul	99
9) Fluorene	15.31	166	6720	0.363	ng/ul	100
11) Pentachlorophenol	16.71	266	380	0.280	ng/ul	98
12) Phenanthrene	17.05	178	10964	0.374	ng/ul	99
13) Anthracene	17.15	178	8834	0.349	ng/ul	99
15) Fluoranthene	19.08	202	13097	0.362	ng/ul	99
17) Pyrene	19.45	202	13963	0.392	ng/ul	100
18) Benzo(a)anthracene	21.22	228	11506	0.360	ng/ul	100
19) Chrysene	21.27	228	14489	0.406	ng/ul	99
21) Benzo(b)fluoranthene	22.80	252	13216	0.355	ng/ul	97
22) Benzo(k)fluoranthene	22.84	252	14522	0.373	ng/ul	98
23) Benzo(a)pyrene	23.37	252	13154	0.361	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.68	276	15383	0.417	ng/ul#	93
25) Dibenzo(a,h)anthracene	25.68	278	11749	0.400	ng/ul	96
26) Benzo(g,h,i)perylene	26.37	276	13693	0.444	ng/ul	97

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

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Instrument :
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ClientSampleId :
SSTD0.427

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