

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN080320\
 Data File : BN011328.D
 Acq On : 03 Aug 2020 09:03
 Operator : CG/JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.411

Quant Time: Aug 03 10:08:09 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SOM-EPA-SIM-BN073020.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Aug 03 10:06:56 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.47	152	1193	0.40	ng/ul	0.00
2) Naphthalene-d8	10.21	136	4356	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.10	164	2582	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.85	188	5753	0.40	ng/ul	0.00
16) Chrysene-d12	21.07	240	5623	0.40	ng/ul	0.00
20) Perylene-d12	23.19	264	5903	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.81	152	4534	0.41	ng/ul	0.00
14) Fluoranthene-d10	18.89	212	11425	0.41	ng/ul	0.00
Target Compounds						Qvalue
3) Naphthalene	10.26	128	5362	0.414	ng/ul	99
5) 2-Methylnaphthalene	11.89	142	3604	0.415	ng/ul	98
7) Acenaphthylene	13.81	152	5201	0.422	ng/ul	100
8) Acenaphthene	14.16	153	4144	0.423	ng/ul	100
9) Fluorene	15.16	166	4772	0.414	ng/ul	99
11) Pentachlorophenol	16.51	266	579	0.411	ng/ul	100
12) Phenanthrene	16.89	178	7962	0.418	ng/ul	99
13) Anthracene	16.99	178	6903	0.415	ng/ul	99
15) Fluoranthene	18.92	202	9694	0.412	ng/ul	98
17) Pyrene	19.29	202	9761	0.382	ng/ul	98
18) Benzo(a)anthracene	21.05	228	9253	0.428	ng/ul	100
19) Chrysene	21.10	228	10590	0.428	ng/ul	99
21) Benzo(b)fluoranthene	22.57	252	10421	0.423	ng/ul	99
22) Benzo(k)fluoranthene	22.61	252	11556	0.418	ng/ul	99
23) Benzo(a)pyrene	23.10	252	8982	0.419	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.26	276	10532	0.468	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.27	278	8314	0.464	ng/ul	98
26) Benzo(g,h,i)perylene	25.88	276	8275	0.407	ng/ul	99

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

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