

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005522.D
 Acq On : 07 May 2019 18:42
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.467

Quant Time: May 07 19:16:54 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN050719.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue May 07 07:04:57 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.60	152	2405	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.36	136	9587	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.23	164	5459	0.40	ng/ul	-0.01
10) Phenanthrene-d10	16.99	188	13403	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	13562	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	12979	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.96	152	6140	0.41	ng/ul	-0.03
14) Fluoranthene-d10	19.03	212	16583	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.41	128	10015	0.395	ng/ul	97
5) 2-Methylnaphthalene	12.04	142	7129	0.412	ng/ul	99
7) Acenaphthylene	13.95	152	10973	0.399	ng/ul	99
8) Acenaphthene	14.30	153	7784	0.400	ng/ul	98
9) Fluorene	15.30	166	8969	0.395	ng/ul	99
11) Pentachlorophenol	16.66	266	1248	0.359	ng/ul	90
12) Phenanthrene	17.04	178	14690	0.402	ng/ul	100
13) Anthracene	17.13	178	14606	0.401	ng/ul	99
15) Fluoranthene	19.06	202	19911	0.394	ng/ul	99
17) Pyrene	19.43	202	20880	0.417	ng/ul	98
18) Benzo(a)anthracene	21.20	228	19208	0.411	ng/ul	100
19) Chrysene	21.26	228	21898	0.381	ng/ul	100
21) Benzo(b)fluoranthene	22.77	252	17284	0.443	ng/ul	96
22) Benzo(k)fluoranthene	22.81	252	20447	0.418	ng/ul	99
23) Benzo(a)pyrene	23.33	252	17679	0.405	ng/ul#	93
24) Indeno(1,2,3-cd)pyrene	25.63	276	16259	0.333	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.64	278	13487	0.334	ng/ul	99
26) Benzo(g,h,i)perylene	26.29	276	13259	0.319	ng/ul	100

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

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