

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN050719\
 Data File : BN005496.D
 Acq On : 06 May 2019 21:38
 Operator : JU/SJ
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.465

Quant Time: May 07 04:11:24 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 03:05:33 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	1544	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	5960	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	3420	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	8851	0.40	ng/ul	0.00
16) Chrysene-d12	21.23	240	9666	0.40	ng/ul	0.00
20) Perylene-d12	23.44	264	11377	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	3673	0.39	ng/ul	0.02
14) Fluoranthene-d10	19.04	212	10766	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	6195	0.393	ng/ul	99
5) 2-Methylnaphthalene	12.07	142	4196	0.390	ng/ul	99
7) Acenaphthylene	13.96	152	6699	0.389	ng/ul	98
8) Acenaphthene	14.31	153	4755	0.390	ng/ul	97
9) Fluorene	15.32	166	5584	0.392	ng/ul	98
11) Pentachlorophenol	16.68	266	791	0.345	ng/ul#	84
12) Phenanthrene	17.04	178	9226	0.383	ng/ul	100
13) Anthracene	17.15	178	9113	0.379	ng/ul	100
15) Fluoranthene	19.07	202	12925	0.387	ng/ul	100
17) Pyrene	19.44	202	13705	0.384	ng/ul	99
18) Benzo(a)anthracene	21.21	228	13007	0.390	ng/ul	99
19) Chrysene	21.27	228	15857	0.387	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	13364	0.391	ng/ul	97
22) Benzo(k)fluoranthene	22.83	252	16952	0.395	ng/ul	100
23) Benzo(a)pyrene	23.35	252	15032	0.393	ng/ul	97
24) Indeno(1,2,3-cd)pyrene	25.65	276	16530	0.387	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.66	278	13661	0.387	ng/ul	98
26) Benzo(g,h,i)perylene	26.31	276	14017	0.384	ng/ul	99

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

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