

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110519\
 Data File : BN008450.D
 Acq On : 05 Nov 2019 17:02
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 11 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SST00.423

Quant Time: Nov 07 11:11:42 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN110519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Nov 07 10:57:49 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.57	152	1234	0.40	ng/ul	0.01
2) Naphthalene-d8	10.32	136	5277	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.20	164	3256	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6097	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	4453	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	3933	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	3266	0.39	ng/ul	0.01
14) Fluoranthene-d10	18.99	212	7588	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6120	0.407	ng/ul	99
5) 2-Methylnaphthalene	12.01	142	3911	0.391	ng/ul	99
7) Acenaphthylene	13.92	152	5635	0.371	ng/ul	100
8) Acenaphthene	14.26	153	4705	0.383	ng/ul	98
9) Fluorene	15.26	166	5028	0.374	ng/ul	99
11) Pentachlorophenol	16.63	266	468	0.353	ng/ul	97
12) Phenanthrene	17.00	178	7610	0.411	ng/ul	99
13) Anthracene	17.09	178	6484	0.399	ng/ul	100
15) Fluoranthene	19.02	202	8841	0.416	ng/ul	100
17) Pyrene	19.38	202	9200	0.423	ng/ul	99
18) Benzo(a)anthracene	21.15	228	5913	0.381	ng/ul	99
19) Chrysene	21.20	228	8437	0.428	ng/ul	99
21) Benzo(b)fluoranthene	22.68	252	5686	0.388	ng/ul	92
22) Benzo(k)fluoranthene	22.73	252	7019	0.421	ng/ul	98
23) Benzo(a)pyrene	23.24	252	5849	0.396	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.48	276	6756	0.403	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.50	278	5324	0.400	ng/ul	97
26) Benzo(g,h,i)perylene	26.12	276	6158	0.410	ng/ul	98

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

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