

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110719\
 Data File : BN008497.D
 Acq On : 07 Nov 2019 22:37
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.405

Quant Time: Nov 08 06:47:33 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 : Fri Nov 08 06:44:20 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1238	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5267	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	2863	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6243	0.40	ng/ul	0.00
16) Chrysene-d12	21.15	240	5367	0.40	ng/ul	0.00
20) Perylene-d12	23.32	264	5880	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.92	152	3128	0.38	ng/ul	0.00
14) Fluoranthene-d10	18.98	212	7469	0.40	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6112	0.408	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	3764	0.377	ng/ul	98
7) Acenaphthylene	13.91	152	5025	0.377	ng/ul	99
8) Acenaphthene	14.25	153	4090	0.379	ng/ul	98
9) Fluorene	15.25	166	4589	0.388	ng/ul	100
11) Pentachlorophenol	16.62	266	554	0.408	ng/ul	100
12) Phenanthrene	16.98	178	7641	0.403	ng/ul	99
13) Anthracene	17.08	178	6629	0.399	ng/ul	100
15) Fluoranthene	19.01	202	8844	0.406	ng/ul	98
17) Pyrene	19.38	202	9280	0.354	ng/ul	99
18) Benzo(a)anthracene	21.14	228	7637	0.409	ng/ul	99
19) Chrysene	21.19	228	9586	0.403	ng/ul	99
21) Benzo(b)fluoranthene	22.68	252	8612	0.393	ng/ul	94
22) Benzo(k)fluoranthene	22.72	252	9442	0.379	ng/ul	97
23) Benzo(a)pyrene	23.23	252	8806	0.399	ng/ul	91
24) Indeno(1,2,3-cd)pyrene	25.46	276	10536	0.420	ng/ul#	100
25) Dibenzo(a,h)anthracene	25.47	278	8310	0.418	ng/ul	96
26) Benzo(g,h,i)perylene	26.11	276	9466	0.421	ng/ul	99

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

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