

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN110619\
 Data File : BN008465.D
 Acq On : 06 Nov 2019 09:59
 Operator : JU
 Sample : SSTDC00.4
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.425

Quant Time: Nov 07 12:30:14 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 12:16:36 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.56	152	1248	0.40	ng/ul	0.00
2) Naphthalene-d8	10.32	136	5445	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.19	164	3437	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.95	188	6585	0.40	ng/ul	0.00
16) Chrysene-d12	21.16	240	4996	0.40	ng/ul	0.00
20) Perylene-d12	23.33	264	4411	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.93	152	3347	0.39	ng/ul	0.01
14) Fluoranthene-d10	18.99	212	8332	0.42	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.37	128	6337	0.409	ng/ul	99
5) 2-Methylnaphthalene	12.00	142	4021	0.390	ng/ul	99
7) Acenaphthylene	13.91	152	5948	0.371	ng/ul	100
8) Acenaphthene	14.26	153	5011	0.386	ng/ul	100
9) Fluorene	15.26	166	5340	0.376	ng/ul	99
11) Pentachlorophenol	16.63	266	443	0.309	ng/ul	98
12) Phenanthrene	16.99	178	8201	0.410	ng/ul	100
13) Anthracene	17.09	178	6883	0.393	ng/ul	100
15) Fluoranthene	19.02	202	9780	0.426	ng/ul	98
17) Pyrene	19.39	202	10319	0.423	ng/ul	99
18) Benzo(a)anthracene	21.15	228	6379	0.367	ng/ul	100
19) Chrysene	21.20	228	9904	0.447	ng/ul	99
21) Benzo(b)fluoranthene	22.69	252	6645	0.404	ng/ul	92
22) Benzo(k)fluoranthene	22.73	252	7871	0.421	ng/ul	99
23) Benzo(a)pyrene	23.24	252	6481	0.392	ng/ul#	83
24) Indeno(1,2,3-cd)pyrene	25.48	276	7276	0.387	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.49	278	5620	0.377	ng/ul	99
26) Benzo(g,h,i)perylene	26.12	276	6782	0.402	ng/ul	100

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

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