

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN111819\
 Data File : BN008622.D
 Acq On : 19 Nov 2019 02:08
 Operator : JU
 Sample : SSTDC00.4EC
 Misc :
 ALS Vial : 2 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SSTD0.410

Quant Time: Nov 19 03:24:05 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN111519.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Nov 19 01:47:58 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	3187	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	11802	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.25	164	6638	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.00	188	13538	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	10973	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	11481	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.98	152	7836	0.40	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	16886	0.41	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	14150	0.398	ng/ul	99
5) 2-Methylnaphthalene	12.06	142	9899	0.398	ng/ul	99
7) Acenaphthylene	13.97	152	14830	0.404	ng/ul	99
8) Acenaphthene	14.32	153	10254	0.376	ng/ul	98
9) Fluorene	15.31	166	11765	0.368	ng/ul	100
11) Pentachlorophenol	16.65	266	1911	0.373	ng/ul	99
12) Phenanthrene	17.03	178	18994	0.416	ng/ul	99
13) Anthracene	17.13	178	18775	0.435	ng/ul	100
15) Fluoranthene	19.06	202	22056	0.393	ng/ul	96
17) Pyrene	19.42	202	22750	0.517	ng/ul	100
18) Benzo(a)anthracene	21.17	228	20304	0.460	ng/ul	96
19) Chrysene	21.22	228	19063	0.440	ng/ul	97
21) Benzo(b)fluoranthene	22.72	252	19015	0.402	ng/ul#	76
22) Benzo(k)fluoranthene	22.76	252	18905	0.403	ng/ul#	70
23) Benzo(a)pyrene	23.27	252	18464	0.415	ng/ul#	75
24) Indeno(1,2,3-cd)pyrene	25.51	276	21667	0.417	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.53	278	17506	0.417	ng/ul#	82
26) Benzo(g,h,i)perylene	26.17	276	18248	0.420	ng/ul#	87

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

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