

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008986.D
 Acq On : 14 Dec 2019 02:46
 Operator : JU
 Sample : PB123046BS
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS046

Quant Time: Dec 16 00:51:11 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 : Mon Dec 16 00:48:46 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2565	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	11098	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	7402	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.92	188	16453	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	13066	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	10670	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	7787	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	19634	0.39	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.36	128	13379	0.400	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	9990	0.428	ng/ul	100
7) Acenaphthylene	13.89	152	14747	0.360	ng/ul	100
8) Acenaphthene	14.24	153	11437	0.377	ng/ul	99
9) Fluorene	15.24	166	13428	0.377	ng/ul	99
11) Pentachlorophenol	16.59	266	1640	0.263	ng/ul	100
12) Phenanthrene	16.97	178	22682	0.409	ng/ul	99
13) Anthracene	17.06	178	20507	0.391	ng/ul	100
15) Fluoranthene	18.99	202	26031	0.382	ng/ul	99
17) Pyrene	19.35	202	26062	0.497	ng/ul	98
18) Benzo(a)anthracene	21.11	228	23743	0.452	ng/ul	99
19) Chrysene	21.16	228	23326	0.452	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	19852	0.451	ng/ul	99
22) Benzo(k)fluoranthene	22.68	252	19623	0.450	ng/ul	99
23) Benzo(a)pyrene	23.18	252	17384	0.421	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.38	276	17835	0.369	ng/ul#	98
25) Dibenzo(a,h)anthracene	25.39	278	14383	0.369	ng/ul	98
26) Benzo(g,h,i)perylene	26.01	276	14809	0.367	ng/ul	98

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

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