

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121319\
 Data File : BN008980.D
 Acq On : 13 Dec 2019 23:10
 Operator : JU
 Sample : PB123045BS
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS045

Quant Time: Dec 13 23:51:56 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 : Fri Dec 13 23:22:26 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.55	152	2283	0.40	ng/ul	0.00
2) Naphthalene-d8	10.31	136	10079	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.18	164	6757	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.93	188	15257	0.40	ng/ul	0.00
16) Chrysene-d12	21.13	240	12533	0.40	ng/ul	0.00
20) Perylene-d12	23.27	264	10609	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.90	152	7133	0.43	ng/ul	0.00
14) Fluoranthene-d10	18.96	212	18285	0.39	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.36	128	12196	0.401	ng/ul	100
5) 2-Methylnaphthalene	11.98	142	9215	0.434	ng/ul	100
7) Acenaphthylene	13.89	152	13766	0.368	ng/ul	99
8) Acenaphthene	14.24	153	10574	0.381	ng/ul	98
9) Fluorene	15.24	166	12415	0.381	ng/ul	99
11) Pentachlorophenol	16.59	266	1632	0.283	ng/ul	98
12) Phenanthrene	16.97	178	21027	0.409	ng/ul	99
13) Anthracene	17.06	178	19191	0.395	ng/ul	99
15) Fluoranthene	18.99	202	24396	0.386	ng/ul	99
17) Pyrene	19.35	202	24468	0.487	ng/ul	99
18) Benzo(a)anthracene	21.11	228	22981	0.456	ng/ul	100
19) Chrysene	21.17	228	22407	0.453	ng/ul	100
21) Benzo(b)fluoranthene	22.64	252	19731	0.451	ng/ul	99
22) Benzo(k)fluoranthene	22.68	252	19400	0.447	ng/ul	99
23) Benzo(a)pyrene	23.18	252	17461	0.425	ng/ul	99
24) Indeno(1,2,3-cd)pyrene	25.38	276	17409	0.363	ng/ul	99
25) Dibenzo(a,h)anthracene	25.39	278	14182	0.366	ng/ul	98
26) Benzo(g,h,i)perylene	26.02	276	14511	0.361	ng/ul	98

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

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