

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN122120\
 Data File : BN013236.D
 Acq On : 21 Dec 2020 11:59
 Operator : CG/JU
 Sample : L5151-06
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AE8

Manual Integrations
 APPROVED

mohammad
 12/22/2020 2:39:40 PM

Quant Time: Dec 21 12:31:37 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN120320.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Dec 21 10:25:14 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.33	152	757	0.40	ng/ul	0.00
2) Naphthalene-d8	10.09	136	2666	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.00	164	1550	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.76	188	3111m	0.40	ng/ul	0.00
16) Chrysene-d12	20.99	240	2277	0.40	ng/ul	0.00
20) Perylene-d12	23.08	264	2417	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.69	152	987	0.25	ng/ul	0.00
14) Fluoranthene-d10	18.81	212	2013	0.24	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.14	128	23727	3.076	ng/ul#	91
5) 2-Methylnaphthalene	11.77	142	18630	3.676	ng/ul	99
7) Acenaphthylene	13.72	152	13260	1.807	ng/ul	96
8) Acenaphthene	14.07	153	4446	0.739	ng/ul	97
9) Fluorene	15.06	166	7273	1.053	ng/ul	99
12) Phenanthrene	16.81	178	135752	13.015	ng/ul	96
13) Anthracene	16.89	178	33214	3.742	ng/ul	95
15) Fluoranthene	18.84	202	312103	25.179	ng/ul	98
17) Pyrene	19.21	202	259258	24.105	ng/ul	99
18) Benzo(a)anthracene	20.98	228	183070	19.433	ng/ul	100
19) Chrysene	21.03	228	173591	16.677	ng/ul	98
21) Benzo(b)fluoranthene	22.47	252	292027m	26.116	ng/ul	
22) Benzo(k)fluoranthene	22.51	252	97885m	7.920	ng/ul	
23) Benzo(a)pyrene	22.99	252	198420m	19.124	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.12	276	136974	10.050	ng/ul#	97
25) Dibenzo(a,h)anthracene	25.12	278	34428	3.194	ng/ul	95
26) Benzo(a,h,i)perylene	25.74	276	107327	8.706	ng/ul	96

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

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