

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN090420\
 Data File : BN011723.D
 Acq On : 05 Sep 2020 05:00
 Operator : CG/JU
 Sample : L3843-02
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AC1

Manual Integrations
 APPROVED

mohammad
 9/8/2020 1:47:39 PM

Quant Time: Sep 05 06:36:40 2020
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN081220.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 05 05:40:34 2020
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.73	152	2	0.40	ng/ul	0.00
2) Naphthalene-d8	10.52	136	81	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.38	164	496	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.12	188	539m	0.40	ng/ul	0.00
16) Chrysene-d12	21.31	240	631	0.40	ng/ul	0.00
20) Perylene-d12	23.56	264	325m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.12	152	5027	43.85	ng/ul	0.00
14) Fluoranthene-d10	19.15	212	10329	6.81	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.57	128	2133	9.034	ng/ul	96
5) 2-Methylnaphthalene	12.19	142	360	2.412	ng/ul	96
7) Acenaphthylene	14.10	152	1865	0.747	ng/ul	95
8) Acenaphthene	14.44	153	418	0.215	ng/ul#	92
9) Fluorene	15.43	166	423	0.185	ng/ul#	89
12) Phenanthrene	17.16	178	1440	0.790	ng/ul#	91
13) Anthracene	17.25	178	1957	1.305	ng/ul#	93
15) Fluoranthene	19.18	202	4457	2.125	ng/ul	91
17) Pyrene	19.55	202	13600m	5.514	ng/ul	
18) Benzo(a)anthracene	21.29	228	3629	1.714	ng/ul	92
19) Chrysene	21.35	228	7871	2.757	ng/ul	97
21) Benzo(b)fluoranthene	22.89	252	14752m	10.443	ng/ul	
22) Benzo(k)fluoranthene	22.93	252	5226m	3.346	ng/ul	
23) Benzo(a)pyrene	23.46	252	8815	7.427	ng/ul#	91
24) Indeno(1,2,3-cd)pyrene	25.82	276	5681	3.540	ng/ul	100
25) Dibenzo(a,h)anthracene	25.82	278	1641	1.254	ng/ul#	59
26) Benzo(a,h,i)perylene	26.51	276	5410	3.601	ng/ul	94

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

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