

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008094.D
 Acq On : 12 Oct 2019 02:14
 Operator : JU
 Sample : K5124-05
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLM67

Manual Integrations
 APPROVED

mohammad
 10/15/2019 7:57:37 AM

Quant Time: Oct 13 01:29:53 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN101019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Oct 11 19:05:54 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	2471	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	11618	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	7548	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	17045m	0.40	ng/ul	0.00
16) Chrysene-d12	21.21	240	17691m	0.40	ng/ul	0.00
20) Perylene-d12	23.40	264	21224	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	3701	0.19	ng/ul	0.00
14) Fluoranthene-d10	19.04	212	11928m	0.20	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	2077	0.063	ng/ul#	88
5) 2-Methylnaphthalene	12.07	142	891	0.039	ng/ul	96
11) Pentachlorophenol	16.68	266	109m	0.026	ng/ul	
12) Phenanthrene	17.05	178	4394m	0.088	ng/ul	
15) Fluoranthene	19.07	202	7043m	0.107	ng/ul	
17) Pyrene	19.44	202	7165m	0.102	ng/ul	
18) Benzo(a)anthracene	21.19	228	3462	0.054	ng/ul#	54
19) Chrysene	21.24	228	7578	0.096	ng/ul#	81
21) Benzo(b)fluoranthene	22.75	252	8400m	0.122	ng/ul	
22) Benzo(k)fluoranthene	22.79	252	2327m	0.030	ng/ul	
23) Benzo(a)pyrene	23.30	252	5945	0.082	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.57	276	2709	0.032	ng/ul#	57
26) Benzo(g,h,i)perylene	26.24	276	8550	0.109	ng/ul#	68

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

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