

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101119\
 Data File : BN008102.D
 Acq On : 12 Oct 2019 08:11
 Operator : JU
 Sample : K5124-16
 Misc :
 ALS Vial : 35 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLM92

Manual Integrations
 APPROVED

mohammad
 10/15/2019 7:58:13 AM

Quant Time: Oct 13 02:12:51 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	3515	0.40	ng/ul	-0.01
2) Naphthalene-d8	10.39	136	15577	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.26	164	11417	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.01	188	24689m	0.40	ng/ul	0.00
16) Chrysene-d12	21.22	240	25689m	0.40	ng/ul	0.00
20) Perylene-d12	23.43	264	17391m	0.40	ng/ul	0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.99	152	6078	0.23	ng/ul	-0.01
14) Fluoranthene-d10	19.05	212	18273m	0.21	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.44	128	129834	2.928	ng/ul	97
5) 2-Methylnaphthalene	12.06	142	78547	2.582	ng/ul	99
7) Acenaphthylene	13.98	152	6491	0.119	ng/ul#	80
8) Acenaphthene	14.32	153	11432	0.270	ng/ul	99
9) Fluorene	15.31	166	12276	0.256	ng/ul#	68
12) Phenanthrene	17.05	178	89112	1.233	ng/ul	96
13) Anthracene	17.14	178	22112m	0.348	ng/ul	
15) Fluoranthene	19.07	202	118591m	1.242	ng/ul	
17) Pyrene	19.44	202	301610	2.949	ng/ul	99
18) Benzo(a)anthracene	21.20	228	63606m	0.681	ng/ul	
19) Chrysene	21.25	228	102039m	0.890	ng/ul	
21) Benzo(b)fluoranthene	22.77	252	111192m	1.964	ng/ul	
22) Benzo(k)fluoranthene	22.80	252	21255m	0.331	ng/ul	
23) Benzo(a)pyrene	23.33	252	51503	0.865	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.61	276	34800	0.496	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.61	278	17805	0.321	ng/ul#	1
26) Benzo(a,h,i)perylene	26.29	276	102360	1.591	ng/ul#	81

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

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