

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008205.D
 Acq On : 16 Oct 2019 19:07
 Operator : JU
 Sample : K5124-02DL 5X
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLM51DL

Manual Integrations
 APPROVED

mohammad
 10/18/2019 9:13:44 AM

Quant Time: Oct 17 12:12:09 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 : Thu Oct 17 01:48:53 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.61	152	3486	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	16490	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	10237	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	20786m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	13485	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	13880	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	1044	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	1999	0.03	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	22950	0.489	ng/ul	98
5) 2-Methylnaphthalene	12.04	142	8959	0.278	ng/ul	100
7) Acenaphthylene	13.96	152	61012	1.251	ng/ul	100
8) Acenaphthene	14.31	153	22075	0.582	ng/ul	100
9) Fluorene	15.30	166	26263	0.610	ng/ul	97
12) Phenanthrene	17.03	178	754436	12.397	ng/ul	99
13) Anthracene	17.12	178	71727	1.339	ng/ul	99
15) Fluoranthene	19.06	202	1025013	12.753	ng/ul	99
17) Pyrene	19.42	202	962005	17.917	ng/ul	99
18) Benzo(a)anthracene	21.18	228	324286	6.614	ng/ul	96
19) Chrysene	21.23	228	401001	6.665	ng/ul	97
21) Benzo(b)fluoranthene	22.73	252	549140m	12.153	ng/ul	
22) Benzo(k)fluoranthene	22.77	252	162788m	3.180	ng/ul	
23) Benzo(a)pyrene	23.29	252	409751	8.618	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.55	276	287926	5.141	ng/ul#	90
25) Dibenzo(a,h)anthracene	25.56	278	60930	1.376	ng/ul#	92
26) Benzo(a,h,i)perylene	26.22	276	293133	5.708	ng/ul	99

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

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