

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008209.D
 Acq On : 16 Oct 2019 21:32
 Operator : JU
 Sample : K5124-19DL 2X
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLMA8DL

Manual Integrations
 APPROVED

mohammad
 10/18/2019 9:13:55 AM

Quant Time: Oct 17 02:51:07 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	3178	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	15156	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	9756	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	20969m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	16446	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	13007	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	1942	0.08	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	5511	0.08	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	1506	0.035	ng/ul#	86
5) 2-Methylnaphthalene	12.04	142	1045	0.035	ng/ul	98
7) Acenaphthylene	13.96	152	4956	0.107	ng/ul	97
9) Fluorene	15.30	166	1147	0.028	ng/ul#	90
12) Phenanthrene	17.03	178	19548	0.318	ng/ul	100
13) Anthracene	17.12	178	2813	0.052	ng/ul#	89
15) Fluoranthene	19.06	202	46899	0.578	ng/ul	98
17) Pyrene	19.42	202	47051	0.719	ng/ul	98
18) Benzo(a)anthracene	21.18	228	16283	0.272	ng/ul#	90
19) Chrysene	21.23	228	27294	0.372	ng/ul#	95
21) Benzo(b)fluoranthene	22.73	252	24433m	0.577	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	10530m	0.220	ng/ul	
23) Benzo(a)pyrene	23.28	252	15082	0.338	ng/ul#	78
24) Indeno(1,2,3-cd)pyrene	25.54	276	12471	0.238	ng/ul#	99
25) Dibenzo(a,h)anthracene	25.55	278	2672	0.064	ng/ul#	14
26) Benzo(a,h,i)perylene	26.20	276	12102	0.251	ng/ul#	86

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

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