

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101619\
 Data File : BN008208.D
 Acq On : 16 Oct 2019 20:56
 Operator : JU
 Sample : K5124-16DL 5X
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 JLM92DL

Manual Integrations
 APPROVED

mohammad
 10/18/2019 9:13:49 AM

Quant Time: Oct 17 02:42:07 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	3102	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	14952	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	9689	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	20575m	0.40	ng/ul	0.00
16) Chrysene-d12	21.20	240	14364	0.40	ng/ul	0.00
20) Perylene-d12	23.38	264	12013	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	1013	0.04	ng/ul	0.00
14) Fluoranthene-d10	19.03	212	2622	0.04	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.42	128	21955	0.516	ng/ul	99
5) 2-Methylnaphthalene	12.04	142	13131	0.450	ng/ul	99
7) Acenaphthylene	13.96	152	1192	0.026	ng/ul#	67
8) Acenaphthene	14.31	153	1534m	0.043	ng/ul	
9) Fluorene	15.30	166	2139	0.052	ng/ul#	63
12) Phenanthrene	17.03	178	13696	0.227	ng/ul#	95
13) Anthracene	17.12	178	3172	0.060	ng/ul#	76
15) Fluoranthene	19.06	202	16474	0.207	ng/ul	92
17) Pyrene	19.42	202	41847	0.732	ng/ul#	95
18) Benzo(a)anthracene	21.18	228	7367	0.141	ng/ul#	42
19) Chrysene	21.23	228	12303	0.192	ng/ul#	67
21) Benzo(b)fluoranthene	22.73	252	14272m	0.365	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	3095m	0.070	ng/ul	
23) Benzo(a)pyrene	23.28	252	5462	0.133	ng/ul#	1
24) Indeno(1,2,3-cd)pyrene	25.55	276	4471	0.092	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.54	278	2242	0.059	ng/ul#	1
26) Benzo(a,h,i)perylene	26.22	276	13344	0.300	ng/ul#	73

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

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