

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN102219\
 Data File : BN008373.D
 Acq On : 24 Oct 2019 05:16
 Operator : JU
 Sample : K5235-12
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPL0

Manual Integrations
 APPROVED

mohammad
 10/24/2019 8:07:30 AM

Quant Time: Oct 24 06:00:08 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 : Wed Oct 23 14:07:55 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.58	152	1893	0.40	ng/ul	-0.02
2) Naphthalene-d8	10.34	136	9279	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.21	164	5728	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.96	188	12400m	0.40	ng/ul	-0.01
16) Chrysene-d12	21.17	240	10355	0.40	ng/ul	-0.01
20) Perylene-d12	23.33	264	9226	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.94	152	3756	0.24	ng/ul	-0.02
14) Fluoranthene-d10	19.00	212	10864	0.25	ng/ul	-0.02
Target Compounds				Ovalue		
12) Phenanthrene	17.00	178	1425	0.039	ng/ul#	89
15) Fluoranthene	19.03	202	3842	0.080	ng/ul	93
17) Pyrene	19.39	202	3396	0.082	ng/ul#	85
18) Benzo(a)anthracene	21.15	228	1685	0.045	ng/ul#	89
19) Chrysene	21.20	228	1792	0.039	ng/ul#	91
21) Benzo(b)fluoranthene	22.69	252	2124m	0.071	ng/ul	
22) Benzo(k)fluoranthene	22.73	252	906m	0.027	ng/ul	
23) Benzo(a)pyrene	23.24	252	1323	0.042	ng/ul#	6
24) Indeno(1,2,3-cd)pyrene	25.49	276	831	0.022	ng/ul#	70
26) Benzo(g,h,i)perylene	26.14	276	801	0.023	ng/ul#	57

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

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