

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN101719\
 Data File : BN008241.D
 Acq On : 18 Oct 2019 03:09
 Operator : JU
 Sample : K5177-15
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 BEPK1

Manual Integrations
 APPROVED

mohammad
 10/20/2019 8:09:58 PM

Quant Time: Oct 18 04:14:22 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	7.60	152	3228	0.40	ng/ul	0.00
2) Naphthalene-d8	10.37	136	15549	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.24	164	9507	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.99	188	20100m	0.40	ng/ul	0.00
16) Chrysene-d12	21.19	240	14394	0.40	ng/ul	0.00
20) Perylene-d12	23.36	264	12774	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.97	152	3320	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.02	212	8401	0.12	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.03	178	5429	0.092	ng/ul	97
15) Fluoranthene	19.05	202	14421	0.186	ng/ul	97
17) Pyrene	19.41	202	12497m	0.218	ng/ul	
18) Benzo(a)anthracene	21.17	228	5428	0.104	ng/ul	92
19) Chrysene	21.22	228	7245	0.113	ng/ul	94
21) Benzo(b)fluoranthene	22.72	252	8812m	0.212	ng/ul	
22) Benzo(k)fluoranthene	22.76	252	2945m	0.063	ng/ul	
23) Benzo(a)pyrene	23.27	252	5238	0.120	ng/ul#	64
24) Indeno(1,2,3-cd)pyrene	25.53	276	4264	0.083	ng/ul#	83
25) Dibenzo(a,h)anthracene	25.53	278	881	0.022	ng/ul#	1
26) Benzo(g,h,i)perylene	26.19	276	4210	0.089	ng/ul#	83

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

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