

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009079.D
 Acq On : 20 Dec 2019 22:14
 Operator : JU
 Sample : K6171-18DL 5X
 Misc :
 ALS Vial : 50 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BZ8DL

Manual Integrations
 APPROVED

mohammad
 12/23/2019 11:55:01 AM

Quant Time: Dec 20 23:47:58 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN121619.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Dec 20 23:31:06 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.54	152	3934	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	14703	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.17	164	7756	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	14544	0.40	ng/ul	0.00
16) Chrysene-d12	21.11	240	8278	0.40	ng/ul	0.00
20) Perylene-d12	23.25	264	8128	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	696	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.94	212	1180	0.03	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.95	178	21505	0.425	ng/ul	100
13) Anthracene	17.04	178	2397	0.053	ng/ul#	93
15) Fluoranthene	18.98	202	54693	0.946	ng/ul	100
17) Pyrene	19.34	202	43755	1.036	ng/ul	99
18) Benzo(a)anthracene	21.10	228	16017	0.436	ng/ul	98
19) Chrysene	21.15	228	20890	0.565	ng/ul	97
21) Benzo(b)fluoranthene	22.62	252	25583m	0.674	ng/ul	
22) Benzo(k)fluoranthene	22.66	252	9163m	0.239	ng/ul	
23) Benzo(a)pyrene	23.16	252	15348	0.461	ng/ul	98
24) Indeno(1,2,3-cd)pyrene	25.35	276	11947	0.330	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.35	278	2685	0.093	ng/ul#	71
26) Benzo(g,h,i)perylene	25.98	276	12015	0.394	ng/ul	98

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

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