

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN121919\
 Data File : BN009078.D
 Acq On : 20 Dec 2019 21:38
 Operator : JU
 Sample : K6171-17DL 5X
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0BZ5DL

Manual Integrations
 APPROVED

mohammad
 12/23/2019 11:54:47 AM

Quant Time: Dec 20 23:45:32 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	3940	0.40	ng/ul	0.00
2) Naphthalene-d8	10.29	136	14951	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.16	164	8030	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.91	188	15698	0.40	ng/ul	0.00
16) Chrysene-d12	21.12	240	8708	0.40	ng/ul	0.00
20) Perylene-d12	23.25	264	7882	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.89	152	670	0.03	ng/ul	0.00
14) Fluoranthene-d10	18.94	212	1247	0.03	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	16.95	178	10308	0.189	ng/ul	99
13) Anthracene	17.04	178	1515	0.031	ng/ul#	89
15) Fluoranthene	18.98	202	27331	0.438	ng/ul	99
17) Pyrene	19.34	202	22469	0.506	ng/ul	98
18) Benzo(a)anthracene	21.10	228	7648	0.198	ng/ul	96
19) Chrysene	21.15	228	9952	0.256	ng/ul	98
21) Benzo(b)fluoranthene	22.62	252	12250m	0.333	ng/ul	
22) Benzo(k)fluoranthene	22.66	252	3878m	0.104	ng/ul	
23) Benzo(a)pyrene	23.16	252	7004	0.217	ng/ul#	87
24) Indeno(1,2,3-cd)pyrene	25.35	276	5489	0.156	ng/ul#	92
25) Dibenzo(a,h)anthracene	25.35	278	1166	0.042	ng/ul#	45
26) Benzo(g,h,i)perylene	25.98	276	5602	0.189	ng/ul	95

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

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