

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007479.D
 Acq On : 28 Aug 2019 21:36
 Operator : HP/JU
 Sample : K4373-20
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AG8

Manual Integrations
 APPROVED

mohammad
 8/29/2019 7:51:02 AM

Quant Time: Aug 28 23:53:43 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN082819.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 20:30:31 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.40	152	5030	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	21882	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	11722	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	27303m	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	22313m	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	19860m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.75	152	10485	0.31	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	26027m	0.29	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	27675	0.448	ng/ul	100
5) 2-Methylnaphthalene	11.83	142	1771	0.043	ng/ul	98
8) Acenaphthene	14.11	153	3941	0.084	ng/ul	98
9) Fluorene	15.11	166	2682	0.053	ng/ul#	97
12) Phenanthrene	16.84	178	14262	0.167	ng/ul	98
13) Anthracene	16.93	178	3020m	0.038	ng/ul	
15) Fluoranthene	18.87	202	14883m	0.131	ng/ul	
17) Pyrene	19.24	202	10702m	0.098	ng/ul	
18) Benzo(a)anthracene	21.00	228	5705	0.057	ng/ul#	90
19) Chrysene	21.05	228	5522	0.052	ng/ul#	92
21) Benzo(b)fluoranthene	22.50	252	6794m	0.078	ng/ul	
23) Benzo(a)pyrene	23.03	252	4215m	0.049	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.18	276	2750m	0.028	ng/ul	
26) Benzo(g,h,i)perylene	25.81	276	2599	0.031	ng/ul#	75

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

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