

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN082819\
 Data File : BN007448.D
 Acq On : 28 Aug 2019 00:53
 Operator : HP/JU
 Sample : K4337-16
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0AG4

Manual Integrations
 APPROVED

mohammad
 8/28/2019 7:48:15 AM

Quant Time: Aug 28 01:43:21 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	5229	0.40	ng/ul	0.00
2) Naphthalene-d8	10.15	136	21790	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.05	164	12270	0.40	ng/ul	0.00
10) Phenanthrene-d10	16.80	188	25726m	0.40	ng/ul	0.00
16) Chrysene-d12	21.02	240	21908m	0.40	ng/ul	0.00
20) Perylene-d12	23.12	264	20028m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	11.76	152	9805	0.29	ng/ul	0.00
14) Fluoranthene-d10	18.84	212	25430m	0.30	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.20	128	1547	0.025	ng/ul#	77
12) Phenanthrene	16.84	178	8217	0.102	ng/ul	96
15) Fluoranthene	18.87	202	13524m	0.126	ng/ul	
17) Pyrene	19.24	202	10179m	0.095	ng/ul	
18) Benzo(a)anthracene	21.00	228	5444	0.056	ng/ul#	94
19) Chrysene	21.05	228	5276m	0.050	ng/ul	
21) Benzo(b)fluoranthene	22.50	252	6741m	0.077	ng/ul	
22) Benzo(k)fluoranthene	22.54	252	2826m	0.031	ng/ul	
23) Benzo(a)pyrene	23.03	252	4219m	0.048	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.18	276	3006m	0.031	ng/ul	
26) Benzo(g,h,i)perylene	25.81	276	2864	0.034	ng/ul#	70

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

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