

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN061819\
 Data File : BN006376.D
 Acq On : 18 Jun 2019 14:05
 Operator : JU/SJ
 Sample : K3335-07
 Misc :
 ALS Vial : 9 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4234

Manual Integrations
 APPROVED

mohammad
 6/21/2019 8:38:57 AM

Quant Time: Jun 18 16:49:21 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA N\METHODS\SOM-EPA-SIM-BN053019.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jun 18 15:57:40 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.64	152	4448	0.40	ng/ul	0.00
2) Naphthalene-d8	10.42	136	17825	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.29	164	9652	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.07	188	275273m	0.40	ng/ul	0.02
16) Chrysene-d12	21.27	240	11597m	0.40	ng/ul	0.00
20) Perylene-d12	23.50	264	10425m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	6400	0.23	ng/ul	0.00
14) Fluoranthene-d10	19.09	212	14466m	0.02	ng/ul	0.00
Target Compounds					Ovalue	
3) Naphthalene	10.47	128	3392	0.069	ng/ul#	92
5) 2-Methylnaphthalene	12.09	142	1167	0.034	ng/ul	100
7) Acenaphthylene	14.01	152	7909	0.147	ng/ul	97
8) Acenaphthene	14.35	153	2887	0.078	ng/ul	99
9) Fluorene	15.35	166	3986	0.092	ng/ul	99
12) Phenanthrene	17.09	178	74428m	0.084	ng/ul	
13) Anthracene	17.18	178	20166m	0.024	ng/ul	
15) Fluoranthene	19.12	202	146685m	0.150	ng/ul	
17) Pyrene	19.49	202	131980	2.161	ng/ul	100
18) Benzo(a)anthracene	21.25	228	55442m	1.081	ng/ul	
19) Chrysene	21.30	228	54270	1.127	ng/ul	99
21) Benzo(b)fluoranthene	22.83	252	71810m	1.652	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	27521m	0.650	ng/ul	
23) Benzo(a)pyrene	23.40	252	46947m	1.160	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.74	276	31051m	0.709	ng/ul	
25) Dibenzo(a,h)anthracene	25.74	278	7461m	0.213	ng/ul	
26) Benzo(a,h,i)perylene	26.42	276	25001m	0.681	ng/ul	

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

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