

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006574.D
 Acq On : 25 Jun 2019 17:25
 Operator : HP/JU
 Sample : K3362-14DL 5X
 Misc :
 ALS Vial : 51 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4203DL

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:22:24 PM

Quant Time: Jun 25 18:38:48 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 25 07:15:41 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	5061	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	17475	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	7760	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	14613	0.40	ng/ul	0.00
16) Chrysene-d12	21.27	240	11491	0.40	ng/ul	0.02
20) Perylene-d12	23.52	264	14219	0.40	ng/ul	0.04
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	1318	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2305	0.06	ng/ul	0.00
Target Compounds				Ovalue		
8) Acenaphthene	14.34	153	1106	0.037	ng/ul	97
9) Fluorene	15.34	166	1473	0.042	ng/ul#	96
12) Phenanthrene	17.08	178	49553	1.056	ng/ul	99
13) Anthracene	17.17	178	8249	0.184	ng/ul	96
15) Fluoranthene	19.11	202	139979	2.702	ng/ul	100
17) Pyrene	19.48	202	113694	1.878	ng/ul	97
18) Benzo(a)anthracene	21.26	228	54560	1.074	ng/ul	99
19) Chrysene	21.31	228	64883	1.360	ng/ul	99
21) Benzo(b)fluoranthene	22.86	252	84681m	1.428	ng/ul	
22) Benzo(k)fluoranthene	22.90	252	33611m	0.582	ng/ul	
23) Benzo(a)pyrene	23.43	252	53984	0.978	ng/ul#	86
24) Indeno(1,2,3-cd)pyrene	25.77	276	40124	0.671	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.77	278	9346	0.196	ng/ul	99
26) Benzo(g,h,i)perylene	26.45	276	32084	0.641	ng/ul	96

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

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