

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006547.D
 Acq On : 25 Jun 2019 00:32
 Operator : HP/JU
 Sample : K3362-09
 Misc :
 ALS Vial : 25 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4200

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:20:54 PM

Quant Time: Jun 25 04:15:48 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 01:51:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.63	152	9572	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	36380	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	19003	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	38952m	0.40	ng/ul	0.00
16) Chrysene-d12	21.25	240	21479m	0.40	ng/ul	-0.01
20) Perylene-d12	23.48	264	19351m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	11940	0.21	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	24093m	0.23	ng/ul	0.00
Target Compounds				Ovalue		
12) Phenanthrene	17.08	178	4118	0.033	ng/ul	94
15) Fluoranthene	19.11	202	10282m	0.074	ng/ul	
17) Pyrene	19.47	202	7871	0.070	ng/ul#	92
18) Benzo(a)anthracene	21.24	228	2777m	0.029	ng/ul	
19) Chrysene	21.29	228	3619m	0.041	ng/ul	
21) Benzo(b)fluoranthene	22.82	252	4319m	0.054	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	1715m	0.022	ng/ul	
23) Benzo(a)pyrene	23.38	252	2522	0.034	ng/ul#	79
24) Indeno(1,2,3-cd)pyrene	25.72	276	1915	0.024	ng/ul	91
26) Benzo(g,h,i)perylene	26.41	276	1749	0.026	ng/ul#	81

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

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