

Data Path : Z:\SVOASRV\HPCHEM1\BNA_N\DATA\BN062419\
 Data File : BN006560.D
 Acq On : 25 Jun 2019 08:08
 Operator : HP/JU
 Sample : K3360-06DL 5X
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A41Y6DL

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:21:37 PM

Quant Time: Jun 25 11:40:45 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	5652	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	20652	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9773	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	17603	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	10535	0.40	ng/ul	0.00
20) Perylene-d12	23.49	264	13419	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	1457	0.05	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	2335	0.05	ng/ul	0.00
Target Compounds				Qvalue		
3) Naphthalene	10.46	128	1187	0.021	ng/ul#	80
8) Acenaphthene	14.34	153	2735	0.073	ng/ul	98
9) Fluorene	15.34	166	2333	0.053	ng/ul#	98
12) Phenanthrene	17.08	178	30520	0.540	ng/ul	99
13) Anthracene	17.17	178	5714	0.106	ng/ul#	94
15) Fluoranthene	19.11	202	42867	0.687	ng/ul	98
17) Pyrene	19.48	202	35134	0.633	ng/ul	98
18) Benzo(a)anthracene	21.24	228	15147	0.325	ng/ul	98
19) Chrysene	21.29	228	17008	0.389	ng/ul	97
21) Benzo(b)fluoranthene	22.82	252	20500m	0.366	ng/ul	
22) Benzo(k)fluoranthene	22.86	252	8119m	0.149	ng/ul	
23) Benzo(a)pyrene	23.39	252	13306	0.255	ng/ul	95
24) Indeno(1,2,3-cd)pyrene	25.72	276	9806	0.174	ng/ul#	86
25) Dibenzo(a,h)anthracene	25.73	278	2446	0.054	ng/ul#	59
26) Benzo(g,h,i)perylene	26.41	276	7439	0.157	ng/ul	94

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

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