

Data Path : Z:\SVOASRV\HPCHEM1\BNA N\DATA\BN062419\
 Data File : BN006575.D
 Acq On : 25 Jun 2019 17:59
 Operator : HP/JU
 Sample : K3362-15DL 2X
 Misc :
 ALS Vial : 52 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleID :
 A4206DL

Manual Integrations
 APPROVED

mohammad
 6/25/2019 10:22:28 PM

Quant Time: Jun 25 18:38:15 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	5302	0.40	ng/ul	0.00
2) Naphthalene-d8	10.41	136	19460	0.40	ng/ul	0.00
6) Acenaphthene-d10	14.28	164	9207	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.04	188	16851	0.40	ng/ul	0.00
16) Chrysene-d12	21.26	240	9871	0.40	ng/ul	0.01
20) Perylene-d12	23.50	264	13226	0.40	ng/ul	0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	3954	0.13	ng/ul	0.00
14) Fluoranthene-d10	19.08	212	5804	0.13	ng/ul	0.00
Target Compounds				Ovalue		
3) Naphthalene	10.46	128	1838	0.034	ng/ul#	87
7) Acenaphthylene	14.00	152	12289	0.240	ng/ul	98
8) Acenaphthene	14.34	153	1011	0.028	ng/ul	96
9) Fluorene	15.34	166	1747	0.042	ng/ul#	94
12) Phenanthrene	17.08	178	25846	0.478	ng/ul	99
13) Anthracene	17.17	178	9877	0.191	ng/ul	97
15) Fluoranthene	19.11	202	51593	0.864	ng/ul	97
17) Pyrene	19.48	202	49557	0.953	ng/ul	99
18) Benzo(a)anthracene	21.24	228	21626	0.495	ng/ul	97
19) Chrysene	21.30	228	25812	0.630	ng/ul	98
21) Benzo(b)fluoranthene	22.83	252	44068m	0.799	ng/ul	
22) Benzo(k)fluoranthene	22.87	252	17476m	0.325	ng/ul	
23) Benzo(a)pyrene	23.40	252	31767	0.618	ng/ul#	88
24) Indeno(1,2,3-cd)pyrene	25.73	276	27929	0.502	ng/ul#	87
25) Dibenzo(a,h)anthracene	25.73	278	6282	0.141	ng/ul	96
26) Benzo(a,h,i)perylene	26.42	276	25000	0.537	ng/ul	96

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

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