

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021934.D
 Acq On : 08 Aug 2019 15:39
 Operator : HP/JU
 Sample : K4028-18
 Misc :
 ALS Vial : 44 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENR5

Manual Integrations
 APPROVED

Sohil
 8/13/2019 8:29:28 AM

Quant Time: Aug 09 01:12:44 2019
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA M\METHODS\SOM-EPA-SIM-BM080319.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 09 00:56:52 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	826	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	3468	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.27	164	1910	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	4581m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	4388m	0.40	ng/ul	-0.02
20) Perylene-d12	23.44	264	4825m	0.40	ng/ul	-0.02
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	981	0.17	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	2860m	0.19	ng/ul	-0.02
Target Compounds					Ovalue	
3) Naphthalene	10.45	128	1799	0.184	ng/ul	97
5) 2-Methylnaphthalene	12.09	142	1020	0.156	ng/ul	94
7) Acenaphthylene	14.00	152	3803	0.427	ng/ul#	84
8) Acenaphthene	14.34	153	2373	0.326	ng/ul	97
9) Fluorene	15.33	166	2414	0.298	ng/ul	97
12) Phenanthrene	17.07	178	52977m	3.974	ng/ul	
13) Anthracene	17.16	178	12582m	1.107	ng/ul	
15) Fluoranthene	19.09	202	125515m	7.030	ng/ul	
17) Pyrene	19.46	202	155300	8.613	ng/ul	99
18) Benzo(a)anthracene	21.21	228	95779m	6.102	ng/ul	
19) Chrysene	21.27	228	102309	5.068	ng/ul	100
21) Benzo(b)fluoranthene	22.78	252	109531m	6.732	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	42518m	2.309	ng/ul	
23) Benzo(a)pyrene	23.35	252	90584m	5.474	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.66	276	54081m	2.737	ng/ul	
25) Dibenzo(a,h)anthracene	25.65	278	16768m	1.050	ng/ul	
26) Benzo(a,h,i)perylene	26.34	276	56878m	3.198	ng/ul	

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

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