

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021931.D
 Acq On : 08 Aug 2019 13:44
 Operator : HP/JU
 Sample : K4028-14
 Misc :
 ALS Vial : 41 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENQ9

Manual Integrations
 APPROVED

mohammad
 8/9/2019 6:12:43 AM

Quant Time: Aug 08 17:30:51 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 : Thu Aug 08 06:54:24 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	778	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	3418	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.27	164	1895	0.40	ng/ul	-0.01
10) Phenanthrene-d10	17.03	188	4381m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.24	240	4078m	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4055m	0.40	ng/ul	0.00
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.01	152	933	0.17	ng/ul	-0.01
14) Fluoranthene-d10	19.07	212	2611m	0.18	ng/ul	-0.01
Target Compounds				Ovalue		
3) Naphthalene	10.45	128	301	0.031	ng/ul#	57
5) 2-Methylnaphthalene	12.09	142	177	0.028	ng/ul	94
7) Acenaphthylene	14.00	152	672	0.076	ng/ul#	77
8) Acenaphthene	14.34	153	203	0.028	ng/ul#	87
9) Fluorene	15.33	166	223	0.028	ng/ul#	92
12) Phenanthrene	17.07	178	4268	0.335	ng/ul	98
13) Anthracene	17.16	178	1209m	0.111	ng/ul	
15) Fluoranthene	19.10	202	10355m	0.606	ng/ul	
17) Pyrene	19.46	202	11454	0.684	ng/ul	96
18) Benzo(a)anthracene	21.22	228	6580	0.451	ng/ul	97
19) Chrysene	21.27	228	6669	0.355	ng/ul	97
21) Benzo(b)fluoranthene	22.79	252	8591m	0.628	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	3267m	0.211	ng/ul	
23) Benzo(a)pyrene	23.35	252	6721m	0.483	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.67	276	4209m	0.253	ng/ul	
25) Dibenzo(a,h)anthracene	25.68	278	1178m	0.088	ng/ul	
26) Benzo(a,h,i)perylene	26.35	276	4141m	0.277	ng/ul	

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

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