

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021958.D
 Acq On : 09 Aug 2019 07:29
 Operator : HP/JU
 Sample : K4029-05
 Misc :
 ALS Vial : 20 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENX6

Manual Integrations
 APPROVED

Sohil
 8/13/2019 8:30:07 AM

Quant Time: Aug 09 12:41:34 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 01:29:50 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	1066	0.40	ng/ul	0.00
2) Naphthalene-d8	10.39	136	4713	0.40	ng/ul	-0.02
6) Acenaphthene-d10	14.27	164	2602	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	6112m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.23	240	5918m	0.40	ng/ul	-0.01
20) Perylene-d12	23.44	264	6102m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.00	152	1415	0.18	ng/ul	-0.02
14) Fluoranthene-d10	19.06	212	3895m	0.19	ng/ul	-0.02
Target Compounds				Ovalue		
3) Naphthalene	10.44	128	917	0.069	ng/ul#	88
5) 2-Methylnaphthalene	12.08	142	537	0.061	ng/ul	99
7) Acenaphthylene	13.99	152	2722	0.224	ng/ul#	84
8) Acenaphthene	14.33	153	1852	0.187	ng/ul	97
9) Fluorene	15.33	166	1872	0.170	ng/ul	98
12) Phenanthrene	17.07	178	38855m	2.185	ng/ul	
13) Anthracene	17.16	178	8436m	0.556	ng/ul	
15) Fluoranthene	19.09	202	75000m	3.148	ng/ul	
17) Pyrene	19.46	202	73986	3.042	ng/ul	99
18) Benzo(a)anthracene	21.21	228	44144m	2.085	ng/ul	
19) Chrysene	21.27	228	44356	1.629	ng/ul	99
21) Benzo(b)fluoranthene	22.78	252	53347m	2.593	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	20060m	0.861	ng/ul	
23) Benzo(a)pyrene	23.34	252	36447m	1.742	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.65	276	23177m	0.928	ng/ul	
25) Dibenzo(a,h)anthracene	25.65	278	7094m	0.351	ng/ul	
26) Benzo(a,h,i)perylene	26.33	276	18307m	0.814	ng/ul	

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

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