

Data Path : Z:\SVOASRV\HPCHEM1\BNA M\DATA\BM080719\
 Data File : BM021913.D
 Acq On : 08 Aug 2019 01:34
 Operator : HP/JU
 Sample : K4029-06
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 BENX7

Manual Integrations
 APPROVED

mohammad
 8/9/2019 6:11:54 AM

Quant Time: Aug 08 06:45:45 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 04:39:13 2019
 Response via : Initial Calibration

Internal Standards	R.T.	QIon	Response	Conc	Units	Dev(Min)
1) 1,4-Dichlorobenzene-d4	7.62	152	897	0.40	ng/ul	0.00
2) Naphthalene-d8	10.40	136	3925	0.40	ng/ul	-0.01
6) Acenaphthene-d10	14.28	164	2160	0.40	ng/ul	0.00
10) Phenanthrene-d10	17.03	188	5046m	0.40	ng/ul	-0.02
16) Chrysene-d12	21.24	240	4709m	0.40	ng/ul	0.00
20) Perylene-d12	23.45	264	4650m	0.40	ng/ul	-0.01
System Monitoring Compounds						
4) 2-Methylnaphthalene-d10	12.02	152	1110	0.17	ng/ul	0.00
14) Fluoranthene-d10	19.07	212	2999m	0.18	ng/ul	-0.01
Target Compounds				Ovalue		
7) Acenaphthylene	14.00	152	466	0.046	ng/ul#	74
9) Fluorene	15.34	166	245	0.027	ng/ul#	92
12) Phenanthrene	17.07	178	4169	0.284	ng/ul	99
13) Anthracene	17.17	178	1170m	0.093	ng/ul	
15) Fluoranthene	19.10	202	8816m	0.448	ng/ul	
17) Pyrene	19.46	202	10385	0.537	ng/ul	96
18) Benzo(a)anthracene	21.22	228	6139	0.364	ng/ul	97
19) Chrysene	21.27	228	5682	0.262	ng/ul	96
21) Benzo(b)fluoranthene	22.79	252	6686m	0.426	ng/ul	
22) Benzo(k)fluoranthene	22.82	252	2253m	0.127	ng/ul	
23) Benzo(a)pyrene	23.35	252	4956m	0.311	ng/ul	
24) Indeno(1,2,3-cd)pyrene	25.67	276	2893m	0.152	ng/ul	
25) Dibenzo(a,h)anthracene	25.66	278	856m	0.056	ng/ul	
26) Benzo(g,h,i)perylene	26.35	276	2370m	0.138	ng/ul	

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

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