

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016074.D
 Acq On : 26 Aug 2021 16:08
 Operator : CG/JU
 Sample : M3395-17
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 X3625

Quant Time: Aug 26 16:38:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN082621.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Aug 26 11:35:43 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	1235	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	5176	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	3027	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	6425	0.400	ng/ul	0.00
17) Chrysene-d12	21.448	240	6598	0.400	ng/ul	0.02
23) Perylene-d12	23.842	264	5844	0.400	ng/ul	0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.315	96	1040	0.622	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	2702	0.347	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	6343	0.357	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.716	128	3182	0.238	ng/ul	97
7) 2-Methylnaphthalene	12.327	142	2270	0.251	ng/ul	99
10) Acenaphthylene	14.224	152	8052	0.753	ng/ul	98
11) Acenaphthene	14.566	153	2869	0.311	ng/ul	99
12) Fluorene	15.552	166	3614	0.331	ng/ul	98
15) Phenanthrene	17.289	178	11649	0.657	ng/ul	99
16) Anthracene	17.381	178	3458	0.226	ng/ul	98
19) Fluoranthene	19.308	202	7891	0.383	ng/ul	99
20) Pyrene	19.671	202	8791	0.424	ng/ul	99
21) Benzo(a)anthracene	21.430	228	14032	0.683	ng/ul	99
22) Chrysene	21.483	228	8218	0.372	ng/ul	100
24) Benzo(b)fluoranthene	23.114	252	6968	0.320	ng/ul	95
26) Benzo(a)pyrene	23.734	252	9054	0.494	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.313	276	8843	0.364	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.330	278	7876	0.386	ng/ul	99
29) Benzo(g,h,i)perylene	27.071	276	8914	0.426	ng/ul	99

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

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