

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082621\
 Data File : BN016071.D
 Acq On : 26 Aug 2021 13:12
 Operator : CG/JU
 Sample : PB138487BS
 Misc :
 ALS Vial : 4 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS487

Quant Time: Aug 26 13:44:53 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	1381	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	5938	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.502	164	3552	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.246	188	7733	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	7905	0.400	ng/ul	0.01
23) Perylene-d12	23.833	264	6841	0.400	ng/ul	0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.319	96	1123	0.600	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.256	152	2957	0.331	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	7519	0.353	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.353	88	2867	1.517	ng/ul#	86
5) Naphthalene	10.716	128	4806	0.313	ng/ul	98
7) 2-Methylnaphthalene	12.327	142	3263	0.315	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	3066	0.302	ng/ul	98
10) Acenaphthylene	14.219	152	3976	0.317	ng/ul	99
11) Acenaphthene	14.562	153	3383	0.313	ng/ul	97
12) Fluorene	15.552	166	3966	0.309	ng/ul	100
14) Pentachlorophenol	16.900	266	759	0.469	ng/ul	98
15) Phenanthrene	17.289	178	6785	0.318	ng/ul	99
16) Anthracene	17.381	178	5848	0.318	ng/ul	100
19) Fluoranthene	19.308	202	8054	0.326	ng/ul	99
20) Pyrene	19.671	202	8153	0.328	ng/ul	99
21) Benzo(a)anthracene	21.427	228	8103	0.329	ng/ul	99
22) Chrysene	21.477	228	9107	0.344	ng/ul	99
24) Benzo(b)fluoranthene	23.105	252	8857	0.347	ng/ul	99
25) Benzo(k)fluoranthene	23.155	252	8901	0.346	ng/ul	99
26) Benzo(a)pyrene	23.728	252	7241	0.337	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.307	276	9802	0.345	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.320	278	8270	0.346	ng/ul	98
29) Benzo(g,h,i)perylene	27.058	276	8360	0.341	ng/ul	99

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

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