

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020824\
 Data File : BN029591.D
 Acq On : 08 Feb 2024 23:47
 Operator : MA/JU
 Sample : P1303-22
 Misc :
 ALS Vial : 19 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 C0FK3

Manual Integrations
 APPROVED

Reviewed By :Jagrut Upadhyay 02/09/2024
 Supervised By :mohammad ahmed 02/09/2024

Quant Time: Feb 09 00:34:12 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 02:17:13 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.884	152	6137	0.400	ng/ul	0.00
4) Naphthalene-d8	10.683	136	17003	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.525	164	7644	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.276	188	14073	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	9900	0.400	ng/ul	0.00
23) Perylene-d12	23.843	264	12405	0.400	ng/ul	-0.02
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.332	96	44291	5.788	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.278	152	7132	0.330	ng/ul	0.00
18) Fluoranthene-d10	19.309	212	10312	0.317	ng/ul	0.00
Target Compounds						
					Qvalue	
15) Phenanthrene	17.318	178	6110	0.139	ng/ul	97
16) Anthracene	17.411	178	1329	0.036	ng/ul#	84
19) Fluoranthene	19.337	202	21770	0.479	ng/ul	98
20) Pyrene	19.704	202	21091m	0.453	ng/ul	
21) Benzo(a)anthracene	21.458	228	10292	0.288	ng/ul	98
22) Chrysene	21.508	228	8841	0.226	ng/ul	98
24) Benzo(b)fluoranthene	23.124	252	12541m	0.269	ng/ul	
25) Benzo(k)fluoranthene	23.168	252	5553m	0.107	ng/ul	
26) Benzo(a)pyrene	23.741	252	8168	0.186	ng/ul	92
27) Indeno(1,2,3-cd)pyrene	26.295	276	4699	0.088	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.312	278	1144	0.027	ng/ul#	28
29) Benzo(g,h,i)perylene	27.043	276	4660	0.096	ng/ul	93

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

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