

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070223\
 Data File : BN026217.D
 Acq On : 03 Jul 2023 09:25
 Operator : MA/JU
 Sample : 03244-20
 Misc :
 ALS Vial : 74 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 DCGF9

Manual Integrations
 APPROVED

Reviewed By :Yogesh Patel 07/04/2023
 Supervised By :Sohil Jodhani 07/04/2023

Quant Time: Jul 03 23:20:11 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN062223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Jul 01 01:46:24 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.951	152	8059	0.400	ng/ul	-0.01
4) Naphthalene-d8	10.765	136	25580	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.593	164	14427	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.343	188	27437	0.400	ng/ul	0.00
17) Chrysene-d12	21.541	240	15063	0.400	ng/ul	# 0.00
23) Perylene-d12	24.008	264	15709	0.400	ng/ul	0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.318	96	6328	0.697	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.354	152	1949	0.050	ng/ul	-0.01
18) Fluoranthene-d10	19.373	212	3323	0.058	ng/ul	0.00
Target Compounds						
					Qvalue	
10) Acenaphthylene	14.316	152	1987	0.029	ng/ul#	83
15) Phenanthrene	17.385	178	6660	0.073	ng/ul#	93
16) Anthracene	17.478	178	2106	0.026	ng/ul#	75
19) Fluoranthene	19.405	202	19890	0.247	ng/ul	97
20) Pyrene	19.768	202	17121	0.205	ng/ul	99
21) Benzo(a)anthracene	21.523	228	9074	0.146	ng/ul#	81
22) Chrysene	21.576	228	11391	0.176	ng/ul#	85
24) Benzo(b)fluoranthene	23.253	252	18105m	0.270	ng/ul	
25) Benzo(k)fluoranthene	23.294	252	7344m	0.109	ng/ul	
26) Benzo(a)pyrene	23.902	252	11436	0.196	ng/ul#	69
27) Indeno(1,2,3-cd)pyrene	26.591	276	9912	0.156	ng/ul#	92
28) Dibenzo(a,h)anthracene	26.594	278	2386	0.050	ng/ul#	1
29) Benzo(g,h,i)perylene	27.399	276	10559	0.187	ng/ul#	85

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

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