

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070124\
 Data File : BN032361.D
 Acq On : 01 Jul 2024 14:00
 Operator : MA/JU
 Sample : P2871-21DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DN7DL

Quant Time: Jul 01 14:26:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN062724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jun 27 13:23:38 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	6453	0.400	ng/ul	0.00
4) Naphthalene-d8	10.387	136	19424	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.258	164	10613	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.016	188	18742	0.400	ng/ul	0.00
17) Chrysene-d12	21.220	240	16259	0.400	ng/ul	0.00
23) Perylene-d12	23.438	264	19518	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.156	96	426	0.059	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	814	0.028	ng/ul	0.00
18) Fluoranthene-d10	19.050	212	1553	0.032	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.437	128	42402	0.825	ng/ul	100
10) Acenaphthylene	13.980	152	13464	0.305	ng/ul	99
11) Acenaphthene	14.322	153	27791	0.811	ng/ul	99
12) Fluorene	15.317	166	12676	0.311	ng/ul	99
15) Phenanthrene	17.054	178	14202	0.276	ng/ul	99
16) Anthracene	17.147	178	65985	1.440	ng/ul	100
19) Fluoranthene	19.082	202	23414	0.378	ng/ul	98
20) Pyrene	19.449	202	11660	0.180	ng/ul	99
21) Benzo(a)anthracene	21.205	228	22450	0.383	ng/ul	98
22) Chrysene	21.258	228	10265	0.159	ng/ul	95
24) Benzo(b)fluoranthene	22.772	252	21716	0.256	ng/ul	87
25) Benzo(k)fluoranthene	22.816	252	27177	0.330	ng/ul#	90
26) Benzo(a)pyrene	23.345	252	8282	0.140	ng/ul#	54
27) Indeno(1,2,3-cd)pyrene	25.676	276	50724	0.549	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.686	278	36249	0.528	ng/ul	97
29) Benzo(g,h,i)perylene	26.360	276	31356	0.426	ng/ul	97

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

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