

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN070524\
 Data File : BN032506.D
 Acq On : 05 Jul 2024 12:35
 Operator : MA/JU
 Sample : P3010-04DL 10X
 Misc :
 ALS Vial : 7 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 A4DP7DL

Quant Time: Jul 05 13:01:06 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 : Wed Jul 03 03:26:42 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.599	152	4529	0.400	ng/ul	0.00
4) Naphthalene-d8	10.371	136	13495	0.400	ng/ul	#-0.01
9) Acenaphthene-d10	14.244	164	7393	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.003	188	13819	0.400	ng/ul	# 0.00
17) Chrysene-d12	21.214	240	12730	0.400	ng/ul	0.00
23) Perylene-d12	23.430	264	15618	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	256	0.051	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.988	152	442	0.022	ng/ul	0.01
18) Fluoranthene-d10	19.041	212	1095	0.029	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.420	128	51431	1.440	ng/ul	100
10) Acenaphthylene	13.966	152	40586	1.320	ng/ul	99
11) Acenaphthene	14.309	153	33993	1.424	ng/ul	99
12) Fluorene	15.308	166	9394	0.330	ng/ul	100
15) Phenanthrene	17.050	178	21223	0.560	ng/ul	99
16) Anthracene	17.143	178	26310	0.779	ng/ul	99
19) Fluoranthene	19.073	202	43989	0.907	ng/ul	99
20) Pyrene	19.440	202	14430	0.285	ng/ul	99
21) Benzo(a)anthracene	21.197	228	27974	0.609	ng/ul	98
22) Chrysene	21.249	228	12341	0.245	ng/ul	94
24) Benzo(b)fluoranthene	22.763	252	35317	0.520	ng/ul	96
25) Benzo(k)fluoranthene	22.807	252	32396	0.491	ng/ul#	95
26) Benzo(a)pyrene	23.333	252	16889	0.358	ng/ul#	85
27) Indeno(1,2,3-cd)pyrene	25.670	276	20117	0.272	ng/ul	94
28) Dibenzo(a,h)anthracene	25.680	278	5564	0.101	ng/ul#	51
29) Benzo(g,h,i)perylene	26.354	276	35550	0.604	ng/ul	97

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

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