

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM071524\
 Data File : BM046594.D
 Acq On : 15 Jul 2024 12:36
 Operator : MA/JU
 Sample : P3101-01DL 10X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 A4DP2DL

Quant Time: Jul 15 13:16:01 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM071024.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Jul 15 08:57:27 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.506	152	2726	0.400	ng/ul	0.00
4) Naphthalene-d8	10.266	136	7395	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.151	164	5007	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.905	188	11429	0.400	ng/ul	0.00
17) Chrysene-d12	21.114	240	9183	0.400	ng/ul	0.00
23) Perylene-d12	23.253	264	10124	0.400	ng/ul	# 0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.126	96	79	0.035	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.866	152	266	0.022	ng/ul	0.00
18) Fluoranthene-d10	18.942	212	788	0.028	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.315	128	22518	1.197	ng/ul	96
10) Acenaphthylene	13.868	152	23743	1.042	ng/ul	96
11) Acenaphthene	14.215	153	15079	0.987	ng/ul	97
12) Fluorene	15.210	166	10697	0.566	ng/ul	98
15) Phenanthrene	16.943	178	18982	0.582	ng/ul	99
16) Anthracene	17.036	178	12665	0.398	ng/ul	98
19) Fluoranthene	18.974	202	26130	0.693	ng/ul	99
20) Pyrene	19.337	202	31561	0.773	ng/ul	99
21) Benzo(a)anthracene	21.096	228	8545	0.210	ng/ul	99
22) Chrysene	21.146	228	15196	0.387	ng/ul	99
24) Benzo(b)fluoranthene	22.618	252	18481	0.457	ng/ul	97
25) Benzo(k)fluoranthene	22.659	252	14784	0.372	ng/ul	97
26) Benzo(a)pyrene	23.162	252	18098	0.544	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	25.376	276	23120	0.456	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.393	278	15131	0.377	ng/ul	97
29) Benzo(g,h,i)perylene	26.017	276	15802	0.391	ng/ul	99

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

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