

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM102422\
 Data File : BM037200.D
 Acq On : 24 Oct 2022 13:04
 Operator : CG/JU
 Sample : N5103-04
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 COBX1

Quant Time: Oct 25 02:22:42 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM101922.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Oct 22 03:34:37 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.900	152	5369	0.400	ng/ul	0.00
4) Naphthalene-d8	10.699	136	16642	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.530	164	9830	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.272	188	20762	0.400	ng/ul	0.00
17) Chrysene-d12	21.446	240	16657	0.400	ng/ul	0.00
23) Perylene-d12	23.820	264	13862	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	18392	2.810	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.288	152	6183	0.261	ng/ul	0.00
18) Fluoranthene-d10	19.295	212	17294	0.343	ng/ul	0.00
Target Compounds						
					Qvalue	
21) Benzo(a)anthracene	21.432	228	1906	0.029	ng/ul#	89
22) Chrysene	21.484	228	1967	0.026	ng/ul#	91
24) Benzo(b)fluoranthene	23.095	252	1962	0.027	ng/ul#	10
25) Benzo(k)fluoranthene	23.141	252	1814	0.026	ng/ul#	16
26) Benzo(a)pyrene	23.714	252	1539	0.026	ng/ul#	5
27) Indeno(1,2,3-cd)pyrene	26.275	276	1842	0.022	ng/ul#	96
29) Benzo(g,h,i)perylene	27.029	276	1584	0.022	ng/ul#	68

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

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