

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012224\
 Data File : BM043884.D
 Acq On : 22 Jan 2024 13:10
 Operator : MA/JU
 Sample : P1160-08DL 5X
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 XA317DL

Quant Time: Jan 22 13:43:06 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM011924.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Jan 19 22:45:24 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.930	152	5652	0.400	ng/ul	0.00
4) Naphthalene-d8	10.743	136	14721	0.400	ng/ul	-0.02
9) Acenaphthene-d10	14.575	164	7093	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.330	188	14363	0.400	ng/ul	-0.02
17) Chrysene-d12	21.523	240	9962	0.400	ng/ul	-0.01
23) Perylene-d12	23.940	264	12504	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.331	96	770	0.100	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.365	152	1126	0.060	ng/ul	0.00
18) Fluoranthene-d10	19.359	212	2075	0.063	ng/ul	0.00
Target Compounds						
					Qvalue	
5) Naphthalene	10.792	128	69578	1.623	ng/ul	97
10) Acenaphthylene	14.297	152	23072	0.608	ng/ul	99
11) Acenaphthene	14.635	153	47398	1.680	ng/ul	99
12) Fluorene	15.639	166	21526	0.738	ng/ul	99
15) Phenanthrene	17.372	178	25025	0.550	ng/ul	99
16) Anthracene	17.465	178	124145	2.927	ng/ul	97
19) Fluoranthene	19.391	202	39466	0.745	ng/ul	100
20) Pyrene	19.758	202	19302	0.348	ng/ul	99
21) Benzo(a)anthracene	21.509	228	33428	0.863	ng/ul	99
22) Chrysene	21.561	228	16219	0.311	ng/ul	98
24) Benzo(b)fluoranthene	23.201	252	33299	0.700	ng/ul	93
25) Benzo(k)fluoranthene	23.248	252	43867	0.768	ng/ul	90
26) Benzo(a)pyrene	23.838	252	12815	0.264	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	26.440	276	85278	1.361	ng/ul#	98
28) Dibenzo(a,h)anthracene	26.457	278	59793	1.249	ng/ul#	89
29) Benzo(g,h,i)perylene	27.208	276	53190	0.951	ng/ul	96

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

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