

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM012224\
 Data File : BM043885.D
 Acq On : 22 Jan 2024 14:17
 Operator : MA/JU
 Sample : P1160-08DL2 20X
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 XA317DL2

Quant Time: Jan 22 14:49:20 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	7.934	152	5182	0.400 ng/ul	0.00
4) Naphthalene-d8	10.769	136	13421	0.400 ng/ul	0.00
9) Acenaphthene-d10	14.578	164	6656	0.400 ng/ul	0.00
13) Phenanthrene-d10	17.341	188	12778	0.400 ng/ul	0.00
17) Chrysene-d12	21.533	240	8501	0.400 ng/ul	0.00
23) Perylene-d12	23.942	264	10645	0.400 ng/ul	-0.01
System Monitoring Compounds					
3) 1,4-Dioxane-d8	3.335	96	204	0.029 ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.391	152	337	0.020 ng/ul	0.03
18) Fluoranthene-d10	19.371	212	673	0.024 ng/ul	0.00
Target Compounds					
					Qvalue
5) Naphthalene	10.819	128	17128	0.438 ng/ul	99
10) Acenaphthylene	14.315	152	5873	0.165 ng/ul	96
11) Acenaphthene	14.648	153	11881	0.449 ng/ul	98
12) Fluorene	15.661	166	5361	0.196 ng/ul	97
15) Phenanthrene	17.384	178	5947	0.147 ng/ul	95
16) Anthracene	17.481	178	28133	0.746 ng/ul	99
19) Fluoranthene	19.404	202	9528	0.211 ng/ul	97
20) Pyrene	19.766	202	4733	0.100 ng/ul	95
21) Benzo(a)anthracene	21.521	228	7384	0.224 ng/ul	99
22) Chrysene	21.568	228	4287	0.096 ng/ul	98
24) Benzo(b)fluoranthene	23.211	252	7586	0.187 ng/ul	81
25) Benzo(k)fluoranthene	23.258	252	10873	0.224 ng/ul#	89
26) Benzo(a)pyrene	23.845	252	2946	0.071 ng/ul#	28
27) Indeno(1,2,3-cd)pyrene	26.455	276	19731	0.370 ng/ul#	100
28) Dibenzo(a,h)anthracene	26.475	278	13710	0.336 ng/ul	97
29) Benzo(g,h,i)perylene	27.226	276	12745	0.268 ng/ul	99

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

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