

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092223\
 Data File : BM042019.D
 Acq On : 23 Sep 2023 11:13
 Operator : MA/JU
 Sample : PB155517BS
 Misc :
 ALS Vial : 36 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS517

Quant Time: Sep 25 01:04:29 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 13:50:58 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.954	152	5691	0.400	ng/ul	0.00
4) Naphthalene-d8	10.763	136	13058	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.597	164	5662	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.350	188	10680	0.400	ng/ul	0.00
17) Chrysene-d12	21.521	240	4717	0.400	ng/ul	0.00
23) Perylene-d12	23.927	264	4922	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.372	96	5967	0.769	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.346	152	6854	0.419	ng/ul	0.00
18) Fluoranthene-d10	19.371	212	8705	0.448	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.406	88	14355	1.759	ng/ul#	87
5) Naphthalene	10.818	128	16311	0.373	ng/ul	100
7) 2-Methylnaphthalene	12.423	142	9560	0.367	ng/ul	99
8) 1-Methylnaphthalene	12.643	142	10176	0.384	ng/ul	100
10) Acenaphthylene	14.324	152	13875	0.366	ng/ul	100
11) Acenaphthene	14.661	153	10736	0.369	ng/ul	99
12) Fluorene	15.647	166	11822	0.361	ng/ul	100
14) Pentachlorophenol	16.978	266	3842	0.680	ng/ul	99
15) Phenanthrene	17.392	178	18088	0.365	ng/ul	99
16) Anthracene	17.485	178	15990	0.374	ng/ul	100
19) Fluoranthene	19.399	202	19606	0.368	ng/ul	97
20) Pyrene	19.766	202	20175	0.385	ng/ul	99
21) Benzo(a)anthracene	21.504	228	13639	0.381	ng/ul	99
22) Chrysene	21.559	228	14080	0.381	ng/ul	99
24) Benzo(b)fluoranthene	23.187	252	14202	0.375	ng/ul	97
25) Benzo(k)fluoranthene	23.237	252	13998	0.380	ng/ul	97
26) Benzo(a)pyrene	23.822	252	12053	0.387	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	26.424	276	16821	0.393	ng/ul	99
28) Dibenzo(a,h)anthracene	26.441	278	12000	0.388	ng/ul	98
29) Benzo(g,h,i)perylene	27.199	276	14448	0.402	ng/ul	98

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM092223\
Data File : BM042019.D
Acq On : 23 Sep 2023 11:13
Operator : MA/JU
Sample : PB155517BS
Misc :
ALS Vial : 36 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SLCS517

Quant Time: Sep 25 01:04:29 2023
Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\SFAM-EPA-SIM-BM092223.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Sep 22 13:50:58 2023
Response via : Initial Calibration

