

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082622\
 Data File : BM036444.D
 Acq On : 27 Aug 2022 02:59
 Operator : CG/JU
 Sample : PB147281BS
 Misc :
 ALS Vial : 27 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 SLCS281

Quant Time: Aug 27 03:28:49 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM082622.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Aug 26 14:26:34 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.799	152	3195	0.400	ng/ul	0.00
4) Naphthalene-d8	10.600	136	13169	0.400	ng/ul	# 0.00
9) Acenaphthene-d10	14.437	164	7039	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.187	188	14804	0.400	ng/ul	0.00
17) Chrysene-d12	21.365	240	15120	0.400	ng/ul	0.00
23) Perylene-d12	23.691	264	12371	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.205	96	4303	1.000	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.200	152	8591	0.408	ng/ul	0.01
18) Fluoranthene-d10	19.211	212	18342	0.982	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.239	88	10594	2.607	ng/ul#	78
5) Naphthalene	10.649	128	15174	0.409	ng/ul	98
7) 2-Methylnaphthalene	12.272	142	9228	0.414	ng/ul	95
8) 1-Methylnaphthalene	12.481	142	9456	0.419	ng/ul	99
10) Acenaphthylene	14.160	152	13019	0.447	ng/ul	100
11) Acenaphthene	14.497	153	10367	0.441	ng/ul	99
12) Fluorene	15.496	166	11769	0.427	ng/ul	100
14) Pentachlorophenol	16.854	266	2200	0.752	ng/ul	97
15) Phenanthrene	17.225	178	19246	0.364	ng/ul	99
16) Anthracene	17.323	178	16816	0.400	ng/ul	99
19) Fluoranthene	19.244	202	22632	0.428	ng/ul	98
20) Pyrene	19.601	202	23560	0.429	ng/ul	100
21) Benzo(a)anthracene	21.350	228	19326	0.309	ng/ul	99
22) Chrysene	21.400	228	24009	0.279	ng/ul	98
24) Benzo(b)fluoranthene	22.978	252	27836	0.309	ng/ul	90
25) Benzo(k)fluoranthene	23.027	252	26092	0.288	ng/ul#	89
26) Benzo(a)pyrene	23.589	252	21532	0.378	ng/ul#	86
27) Indeno(1,2,3-cd)pyrene	26.104	276	25229	0.586	ng/ul	97
28) Dibenzo(a,h)anthracene	26.121	278	20283	0.556	ng/ul	91
29) Benzo(g,h,i)perylene	26.838	276	22994	0.590	ng/ul	96

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

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM082622\
Data File : BM036444.D
Acq On : 27 Aug 2022 02:59
Operator : CG/JU
Sample : PB147281BS
Misc :
ALS Vial : 27 Sample Multiplier: 1

Instrument :
BNA_M
ClientSampleId :
SLCS281

Quant Time: Aug 27 03:28:49 2022
Quant Method : Z:\SVOASRV\HPCHEM1\BNA_M\METHODS\SFAM-EPA-SIM-BM082622.M
Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
QLast Update : Fri Aug 26 14:26:34 2022
Response via : Initial Calibration

