

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092424\
 Data File : BN034184.D
 Acq On : 24 Sep 2024 03:05
 Operator : RC/JU
 Sample : PB163442BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS442

Quant Time: Sep 24 03:48:42 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Sep 23 19:39:10 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.423	152	6259	0.400	ng/ul	0.00
4) Naphthalene-d8	10.180	136	17964	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.074	164	10211	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.827	188	19187	0.400	ng/ul	0.00
17) Chrysene-d12	21.046	240	11070	0.400	ng/ul	0.00
23) Perylene-d12	23.165	264	10619	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.039	96	5101	0.793	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.780	152	9171	0.356	ng/ul	0.00
18) Fluoranthene-d10	18.870	212	14897	0.451	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.069	88	14849	2.080	ng/ul#	78
5) Naphthalene	10.229	128	17566	0.366	ng/ul	100
7) 2-Methylnaphthalene	11.857	142	11245	0.369	ng/ul	100
8) 1-Methylnaphthalene	12.077	142	13538	0.432	ng/ul	99
10) Acenaphthylene	13.787	152	14779	0.319	ng/ul	100
11) Acenaphthene	14.134	153	12550	0.375	ng/ul	98
12) Fluorene	15.133	166	13680	0.366	ng/ul	100
14) Pentachlorophenol	16.485	266	2384	0.395	ng/ul	100
15) Phenanthrene	16.869	178	20329	0.359	ng/ul	99
16) Anthracene	16.962	178	17058	0.335	ng/ul	99
19) Fluoranthene	18.902	202	19569	0.448	ng/ul	97
20) Pyrene	19.265	202	19941	0.430	ng/ul	97
21) Benzo(a)anthracene	21.029	228	14292	0.328	ng/ul	100
22) Chrysene	21.081	228	16897	0.384	ng/ul	100
24) Benzo(b)fluoranthene	22.539	252	14568	0.353	ng/ul	98
25) Benzo(k)fluoranthene	22.583	252	16221	0.393	ng/ul	98
26) Benzo(a)pyrene	23.074	252	13012	0.393	ng/ul	96
27) Indeno(1,2,3-cd)pyrene	25.245	276	17397	0.350	ng/ul#	92
28) Dibenzo(a,h)anthracene	25.259	278	13293	0.344	ng/ul	96
29) Benzo(g,h,i)perylene	25.869	276	14332	0.369	ng/ul	97

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

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