

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092424\
 Data File : BN034195.D
 Acq On : 24 Sep 2024 10:55
 Operator : RC/JU
 Sample : PB163538BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS538

Quant Time: Sep 24 12:29:22 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	5274	0.400	ng/ul	0.00
4) Naphthalene-d8	10.179	136	14414	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.074	164	7696	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.827	188	14942	0.400	ng/ul	0.00
17) Chrysene-d12	21.049	240	11108	0.400	ng/ul	0.00
23) Perylene-d12	23.168	264	11613	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.039	96	4183	0.771	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.780	152	6921	0.335	ng/ul	0.00
18) Fluoranthene-d10	18.870	212	12462	0.376	ng/ul	0.00
Target Compounds					Qvalue	
2) 1,4-Dioxane	3.069	88	11908	1.979	ng/ul#	78
5) Naphthalene	10.229	128	13388	0.347	ng/ul	100
7) 2-Methylnaphthalene	11.857	142	8278	0.339	ng/ul	99
8) 1-Methylnaphthalene	12.077	142	9982	0.397	ng/ul	100
10) Acenaphthylene	13.787	152	10517	0.301	ng/ul	99
11) Acenaphthene	14.134	153	9032	0.358	ng/ul	100
12) Fluorene	15.133	166	9611	0.341	ng/ul	99
14) Pentachlorophenol	16.489	266	1857	0.395	ng/ul	99
15) Phenanthrene	16.869	178	14967	0.340	ng/ul	100
16) Anthracene	16.962	178	12438	0.314	ng/ul	99
19) Fluoranthene	18.902	202	15929	0.363	ng/ul	98
20) Pyrene	19.265	202	16633	0.358	ng/ul	97
21) Benzo(a)anthracene	21.031	228	13169	0.301	ng/ul	99
22) Chrysene	21.084	228	15736	0.356	ng/ul	99
24) Benzo(b)fluoranthene	22.542	252	14577	0.323	ng/ul	97
25) Benzo(k)fluoranthene	22.583	252	16263	0.361	ng/ul	98
26) Benzo(a)pyrene	23.074	252	13281	0.367	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.245	276	17934	0.330	ng/ul#	91
28) Dibenzo(a,h)anthracene	25.262	278	13691	0.324	ng/ul	97
29) Benzo(g,h,i)perylene	25.872	276	14657	0.345	ng/ul	97

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

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