

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111924\
 Data File : BN035205.D
 Acq On : 20 Nov 2024 18:43
 Operator : RC/JU
 Sample : PB165092BS
 Misc :
 ALS Vial : 18 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS092

Quant Time: Nov 20 23:25:57 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN111624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Nov 18 00:44:37 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.541	152	3536	0.400	ng/ul	0.00
4) Naphthalene-d8	10.300	136	8336	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.180	164	5340	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.937	188	10855	0.400	ng/ul	0.00
17) Chrysene-d12	21.145	240	9424	0.400	ng/ul	0.00
23) Perylene-d12	23.320	264	10371	0.400	ng/ul	#-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.140	96	1904	0.589	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.901	152	4240	0.347	ng/ul	-0.01
18) Fluoranthene-d10	18.977	212	9079	0.349	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.170	88	5215	1.486	ng/ul#	94
5) Naphthalene	10.350	128	6579	0.313	ng/ul	99
7) 2-Methylnaphthalene	11.978	142	4564	0.313	ng/ul	100
8) 1-Methylnaphthalene	12.198	142	4618	0.314	ng/ul	99
10) Acenaphthylene	13.898	152	6584	0.311	ng/ul	100
11) Acenaphthene	14.245	153	5095	0.314	ng/ul	99
12) Fluorene	15.239	166	5787	0.304	ng/ul	98
14) Pentachlorophenol	16.590	266	1528	0.604	ng/ul	98
15) Phenanthrene	16.979	178	9108	0.317	ng/ul	99
16) Anthracene	17.068	178	8245	0.310	ng/ul	100
19) Fluoranthene	19.004	202	10581	0.322	ng/ul	99
20) Pyrene	19.372	202	10885	0.319	ng/ul	99
21) Benzo(a)anthracene	21.131	228	10615	0.325	ng/ul	100
22) Chrysene	21.183	228	11257	0.342	ng/ul	99
24) Benzo(b)fluoranthene	22.673	252	11561	0.337	ng/ul	100
25) Benzo(k)fluoranthene	22.714	252	12177	0.345	ng/ul	99
26) Benzo(a)pyrene	23.223	252	10304	0.329	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	25.476	276	14589	0.345	ng/ul	100
28) Dibenzo(a,h)anthracene	25.490	278	11456	0.345	ng/ul	98
29) Benzo(g,h,i)perylene	26.127	276	11476	0.344	ng/ul	99

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

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