

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033973.D
 Acq On : 16 Sep 2024 04:27
 Operator : JU/RC
 Sample : PB163195BS
 Misc :
 ALS Vial : 65 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS195

Quant Time: Sep 16 06:32:03 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 : Sun Sep 15 01:52:43 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.477	152	4280	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	12014	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.124	164	6061	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.877	188	12510	0.400	ng/ul	0.00
17) Chrysene-d12	21.092	240	10586	0.400	ng/ul	0.00
23) Perylene-d12	23.234	264	11648	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	4063	0.923	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.834	152	7655	0.444	ng/ul	-0.01
18) Fluoranthene-d10	18.920	212	15754	0.499	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.127	88	10402	2.131	ng/ul#	79
5) Naphthalene	10.283	128	12860	0.400	ng/ul	99
7) 2-Methylnaphthalene	11.911	142	7957	0.391	ng/ul	99
8) 1-Methylnaphthalene	12.136	142	9607	0.459	ng/ul	99
10) Acenaphthylene	13.837	152	10243	0.372	ng/ul	99
11) Acenaphthene	14.184	153	8461	0.426	ng/ul	99
12) Fluorene	15.183	166	9045	0.407	ng/ul	99
14) Pentachlorophenol	16.539	266	1165	0.296	ng/ul	99
15) Phenanthrene	16.919	178	14592	0.395	ng/ul	99
16) Anthracene	17.012	178	12695	0.383	ng/ul	100
19) Fluoranthene	18.948	202	17301	0.414	ng/ul	99
20) Pyrene	19.315	202	18237	0.411	ng/ul	99
21) Benzo(a)anthracene	21.077	228	16192	0.389	ng/ul	100
22) Chrysene	21.127	228	18643	0.443	ng/ul	100
24) Benzo(b)fluoranthene	22.599	252	18129	0.401	ng/ul	98
25) Benzo(k)fluoranthene	22.640	252	19675	0.435	ng/ul	98
26) Benzo(a)pyrene	23.140	252	16708	0.460	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.344	276	22436	0.411	ng/ul	99
28) Dibenzo(a,h)anthracene	25.361	278	17366	0.410	ng/ul	97
29) Benzo(g,h,i)perylene	25.985	276	18064	0.424	ng/ul	98

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

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