

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033917.D
 Acq On : 14 Sep 2024 16:24
 Operator : JU/RC
 Sample : PB163391BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS391

Quant Time: Sep 15 01:59:03 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	7.485	152	3702	0.400	ng/ul	0.00
4) Naphthalene-d8	10.243	136	10043	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.131	164	5019	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.884	188	10329	0.400	ng/ul	0.00
17) Chrysene-d12	21.102	240	8952	0.400	ng/ul	0.00
23) Perylene-d12	23.250	264	9833	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.101	96	3037	0.798	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.843	152	5297	0.368	ng/ul	0.00
18) Fluoranthene-d10	18.928	212	10355	0.388	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.131	88	8784	2.080	ng/ul#	79
5) Naphthalene	10.292	128	10319	0.384	ng/ul	100
7) 2-Methylnaphthalene	11.920	142	6394	0.376	ng/ul	99
8) 1-Methylnaphthalene	12.140	142	6435	0.368	ng/ul	99
10) Acenaphthylene	13.845	152	8102	0.355	ng/ul	99
11) Acenaphthene	14.192	153	6657	0.405	ng/ul	99
12) Fluorene	15.191	166	7079	0.385	ng/ul	98
14) Pentachlorophenol	16.542	266	1467	0.451	ng/ul	98
15) Phenanthrene	16.926	178	11114	0.365	ng/ul	99
16) Anthracene	17.019	178	9609	0.351	ng/ul	99
19) Fluoranthene	18.956	202	13137	0.372	ng/ul	99
20) Pyrene	19.323	202	13831	0.369	ng/ul	97
21) Benzo(a)anthracene	21.087	228	12459	0.354	ng/ul	99
22) Chrysene	21.140	228	14353	0.403	ng/ul	100
24) Benzo(b)fluoranthene	22.613	252	13864	0.363	ng/ul	98
25) Benzo(k)fluoranthene	22.654	252	15755	0.413	ng/ul	99
26) Benzo(a)pyrene	23.153	252	13033	0.425	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.363	276	17688	0.384	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.383	278	13696	0.383	ng/ul	96
29) Benzo(g,h,i)perylene	26.010	276	14471	0.402	ng/ul	99

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

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