

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091922\
 Data File : BN021810.D
 Acq On : 19 Sep 2022 18:47
 Operator : CG/JU
 Sample : PB147715BS
 Misc :
 ALS Vial : 15 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS715

Quant Time: Sep 20 00:26:13 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090822.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Sep 20 00:20:21 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.055	152	4501	0.400	ng/ul	0.00
4) Naphthalene-d8	10.883	136	15360	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.712	164	8985	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.464	188	17892	0.400	ng/ul	0.00
17) Chrysene-d12	21.650	240	11077	0.400	ng/ul	# 0.00
23) Perylene-d12	24.158	264	9284	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.368	96	3540	0.639	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.472	152	9163	0.377	ng/ul	0.00
18) Fluoranthene-d10	19.487	212	17863	0.434	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.401	88	9836	1.732	ng/ul#	87
5) Naphthalene	10.938	128	15725	0.363	ng/ul	98
7) 2-Methylnaphthalene	12.544	142	10448	0.371	ng/ul	98
8) 1-Methylnaphthalene	12.764	142	10826	0.375	ng/ul	100
10) Acenaphthylene	14.434	152	15421	0.409	ng/ul	98
11) Acenaphthene	14.772	153	11388	0.356	ng/ul	99
12) Fluorene	15.758	166	13065	0.364	ng/ul	99
14) Pentachlorophenol	17.113	266	2108	0.608	ng/ul	99
15) Phenanthrene	17.506	178	20716	0.367	ng/ul	99
16) Anthracene	17.595	178	18913	0.404	ng/ul	99
19) Fluoranthene	19.520	202	22046	0.426	ng/ul	99
20) Pyrene	19.882	202	21894	0.429	ng/ul	96
21) Benzo(a)anthracene	21.632	228	16443	0.443	ng/ul	99
22) Chrysene	21.688	228	16404	0.405	ng/ul	99
24) Benzo(b)fluoranthene	23.380	252	15169	0.337	ng/ul	89
25) Benzo(k)fluoranthene	23.433	252	14356	0.329	ng/ul#	90
26) Benzo(a)pyrene	24.041	252	14505	0.425	ng/ul#	88
27) Indeno(1,2,3-cd)pyrene	26.797	276	15999	0.417	ng/ul	97
28) Dibenzo(a,h)anthracene	26.821	278	12999	0.399	ng/ul#	92
29) Benzo(g,h,i)perylene	27.616	276	13612	0.385	ng/ul	93

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

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