

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033946.D
 Acq On : 15 Sep 2024 11:39
 Operator : JU/RC
 Sample : PB163355BS
 Misc :
 ALS Vial : 37 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS355

Quant Time: Sep 16 00:13:01 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	7.481	152	3608	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	9965	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.128	164	5065	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.881	188	10660	0.400	ng/ul	0.00
17) Chrysene-d12	21.098	240	9166	0.400	ng/ul	0.00
23) Perylene-d12	23.240	264	9828	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	3671	0.990	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	6671	0.467	ng/ul	0.00
18) Fluoranthene-d10	18.920	212	13851	0.507	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.131	88	9123	2.217	ng/ul#	80
5) Naphthalene	10.288	128	10683	0.401	ng/ul	99
7) 2-Methylnaphthalene	11.916	142	6663	0.395	ng/ul	100
8) 1-Methylnaphthalene	12.136	142	7964	0.459	ng/ul	100
10) Acenaphthylene	13.841	152	8518	0.370	ng/ul	99
11) Acenaphthene	14.188	153	7122	0.429	ng/ul	99
12) Fluorene	15.188	166	7638	0.412	ng/ul	99
14) Pentachlorophenol	16.539	266	902	0.269	ng/ul	98
15) Phenanthrene	16.923	178	12304	0.391	ng/ul	100
16) Anthracene	17.016	178	10512	0.372	ng/ul	99
19) Fluoranthene	18.952	202	14380	0.398	ng/ul	99
20) Pyrene	19.320	202	15156	0.395	ng/ul	97
21) Benzo(a)anthracene	21.083	228	13197	0.366	ng/ul	99
22) Chrysene	21.133	228	15797	0.433	ng/ul	100
24) Benzo(b)fluoranthene	22.605	252	14794	0.388	ng/ul	98
25) Benzo(k)fluoranthene	22.646	252	16937	0.444	ng/ul	99
26) Benzo(a)pyrene	23.146	252	13637	0.445	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.358	276	18468	0.401	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.371	278	14316	0.401	ng/ul	97
29) Benzo(g,h,i)perylene	25.995	276	15175	0.422	ng/ul	99

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

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