

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110321\
 Data File : BN017271.D
 Acq On : 03 Nov 2021 12:10
 Operator : CG/JU
 Sample : PB140312BS
 Misc :
 ALS Vial : 6 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS312

Quant Time: Nov 03 12:39:29 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Nov 03 10:42:03 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.519	152	4010	0.400	ng/ul	0.00
4) Naphthalene-d8	10.289	136	15604	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.171	164	9248	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.924	188	20103	0.400	ng/ul	0.00
17) Chrysene-d12	21.139	240	19399	0.400	ng/ul	0.00
23) Perylene-d12	23.334	264	18377	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.042	96	3281	0.736	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.890	152	8705	0.382	ng/ul	0.00
18) Fluoranthene-d10	18.967	212	21118	0.415	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.072	88	8401	1.754	ng/ul#	80
5) Naphthalene	10.339	128	16292	0.377	ng/ul	100
7) 2-Methylnaphthalene	11.967	142	10816	0.362	ng/ul	99
8) 1-Methylnaphthalene	12.187	142	10877	0.355	ng/ul	100
10) Acenaphthylene	13.889	152	13667	0.374	ng/ul	99
11) Acenaphthene	14.231	153	12543	0.391	ng/ul	99
12) Fluorene	15.230	166	14067	0.375	ng/ul	99
14) Pentachlorophenol	16.586	266	2080	0.459	ng/ul	99
15) Phenanthrene	16.966	178	24304	0.381	ng/ul	100
16) Anthracene	17.059	178	19622	0.371	ng/ul	100
19) Fluoranthene	18.995	202	27995	0.403	ng/ul	99
20) Pyrene	19.357	202	28551	0.401	ng/ul	96
21) Benzo(a)anthracene	21.121	228	22470	0.354	ng/ul	100
22) Chrysene	21.174	228	28680	0.398	ng/ul	99
24) Benzo(b)fluoranthene	22.676	252	27234	0.398	ng/ul	99
25) Benzo(k)fluoranthene	22.720	252	33272	0.428	ng/ul	96
26) Benzo(a)pyrene	23.237	252	23724	0.386	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.543	276	32313	0.371	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.563	278	25303	0.367	ng/ul	98
29) Benzo(g,h,i)perylene	26.210	276	26416	0.358	ng/ul	100

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

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