

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN110321\
 Data File : BN017270.D
 Acq On : 03 Nov 2021 11:34
 Operator : CG/JU
 Sample : PB140310BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS310

Quant Time: Nov 03 12:03:47 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	7.515	152	4124	0.400	ng/ul	0.00
4) Naphthalene-d8	10.289	136	16206	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.171	164	9627	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.928	188	21005	0.400	ng/ul	0.00
17) Chrysene-d12	21.139	240	20439	0.400	ng/ul	0.00
23) Perylene-d12	23.334	264	18665	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.042	96	3290	0.718	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.890	152	8991	0.380	ng/ul	0.00
18) Fluoranthene-d10	18.967	212	22090	0.412	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.072	88	8867	1.800	ng/ul#	78
5) Naphthalene	10.339	128	16895	0.376	ng/ul	100
7) 2-Methylnaphthalene	11.967	142	11318	0.364	ng/ul	100
8) 1-Methylnaphthalene	12.187	142	11246	0.354	ng/ul	100
10) Acenaphthylene	13.884	152	14363	0.377	ng/ul	100
11) Acenaphthene	14.231	153	12998	0.389	ng/ul	99
12) Fluorene	15.230	166	14582	0.374	ng/ul	98
14) Pentachlorophenol	16.586	266	2233	0.472	ng/ul	99
15) Phenanthrene	16.966	178	25453	0.382	ng/ul	99
16) Anthracene	17.059	178	20690	0.374	ng/ul	100
19) Fluoranthene	18.995	202	29091	0.398	ng/ul	100
20) Pyrene	19.362	202	29794	0.397	ng/ul	99
21) Benzo(a)anthracene	21.121	228	23713	0.354	ng/ul	99
22) Chrysene	21.174	228	29835	0.393	ng/ul	99
24) Benzo(b)fluoranthene	22.676	252	28690	0.412	ng/ul	99
25) Benzo(k)fluoranthene	22.720	252	32622	0.414	ng/ul	97
26) Benzo(a)pyrene	23.237	252	24305	0.389	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.543	276	31251	0.353	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.563	278	24637	0.352	ng/ul	99
29) Benzo(g,h,i)perylene	26.214	276	26276	0.351	ng/ul	99

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

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