

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN010322\
 Data File : BN018304.D
 Acq On : 03 Jan 2022 13:04
 Operator : CG/JU
 Sample : PB141604BS
 Misc :
 ALS Vial : 5 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS604

Quant Time: Jan 03 23:59:59 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122721.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Dec 28 00:34:45 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.607	152	1368	0.400	ng/ul	0.00
4) Naphthalene-d8	10.383	136	4836	0.400	ng/ul #	0.00
9) Acenaphthene-d10	14.249	164	3411	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.991	188	8493	0.400	ng/ul	0.00
17) Chrysene-d12	21.189	240	13363	0.400	ng/ul	0.00
23) Perylene-d12	23.410	264	16535	0.400	ng/ul #	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.067	96	1200	0.940	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.989	152	2950	0.351	ng/ul	0.00
18) Fluoranthene-d10	19.023	212	12376	0.313	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.101	88	2441	1.718	ng/ul#	90
5) Naphthalene	10.432	128	4668	0.325	ng/ul	96
7) 2-Methylnaphthalene	12.060	142	3303	0.315	ng/ul	99
8) 1-Methylnaphthalene	12.280	142	3257	0.308	ng/ul	99
10) Acenaphthylene	13.967	152	4601	0.312	ng/ul#	94
11) Acenaphthene	14.310	153	4024	0.339	ng/ul	99
12) Fluorene	15.300	166	5033	0.344	ng/ul	97
14) Pentachlorophenol	16.653	266	1863	0.554	ng/ul	99
15) Phenanthrene	17.033	178	9986	0.355	ng/ul	97
16) Anthracene	17.122	178	8093	0.320	ng/ul	97
19) Fluoranthene	19.055	202	15461	0.291	ng/ul	98
20) Pyrene	19.418	202	17152	0.309	ng/ul	99
21) Benzo(a)anthracene	21.171	228	17362	0.341	ng/ul	99
22) Chrysene	21.224	228	21436	0.388	ng/ul	100
24) Benzo(b)fluoranthene	22.740	252	23967	0.346	ng/ul	95
25) Benzo(k)fluoranthene	22.784	252	28484	0.361	ng/ul	96
26) Benzo(a)pyrene	23.310	252	24443	0.372	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	25.667	276	34526	0.413	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.687	278	26831	0.416	ng/ul	97
29) Benzo(g,h,i)perylene	26.351	276	31644	0.445	ng/ul	99

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

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