

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN12221\
 Data File : BN018169.D
 Acq On : 23 Dec 2021 02:10
 Operator : CG/JU
 Sample : PB141516BS
 Misc :
 ALS Vial : 28 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS516

Quant Time: Dec 23 02:49:55 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN122121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Dec 22 14:08:04 2021
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.611	152	4238	0.400	ng/ul	0.00
4) Naphthalene-d8	10.392	136	13755	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.253	164	8920	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.994	188	20147	0.400	ng/ul	0.00
17) Chrysene-d12	21.190	240	19700	0.400	ng/ul	0.00
23) Perylene-d12	23.414	264	18395	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.076	96	2302	0.526	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.993	152	5209	0.275	ng/ul	0.00
18) Fluoranthene-d10	19.031	212	15397	0.288	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.109	88	6625	1.251	ng/ul	89
5) Naphthalene	10.442	128	10585	0.271	ng/ul	100
7) 2-Methylnaphthalene	12.070	142	6648	0.264	ng/ul	99
8) 1-Methylnaphthalene	12.284	142	6873	0.263	ng/ul	100
10) Acenaphthylene	13.970	152	9462	0.268	ng/ul	99
11) Acenaphthene	14.313	153	8322	0.268	ng/ul	98
12) Fluorene	15.307	166	9528	0.265	ng/ul	98
14) Pentachlorophenol	16.665	266	1300	0.343	ng/ul	97
15) Phenanthrene	17.036	178	17649	0.268	ng/ul	100
16) Anthracene	17.129	178	13774	0.259	ng/ul	100
19) Fluoranthene	19.058	202	21663	0.287	ng/ul	99
20) Pyrene	19.421	202	22682	0.286	ng/ul	99
21) Benzo(a)anthracene	21.176	228	17217	0.263	ng/ul	99
22) Chrysene	21.225	228	23207	0.289	ng/ul	100
24) Benzo(b)fluoranthene	22.745	252	20325	0.289	ng/ul	98
25) Benzo(k)fluoranthene	22.789	252	24894	0.284	ng/ul	100
26) Benzo(a)pyrene	23.318	252	18159	0.272	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	25.675	276	22641	0.258	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.692	278	16999	0.255	ng/ul	98
29) Benzo(g,h,i)perylene	26.363	276	19137	0.264	ng/ul	99

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

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