

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090221\
 Data File : BN016169.D
 Acq On : 03 Sep 2021 11:02
 Operator : CG/JU
 Sample : PB138652BS
 Misc :
 ALS Vial : 39 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS652

Quant Time: Sep 03 11:45:02 2021
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090121.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Sep 02 13:32:21 2021
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.867	152	2655	0.400	ng/ul	0.00
4) Naphthalene-d8	10.666	136	10878	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.501	164	6201	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.250	188	12349	0.400	ng/ul	0.00
17) Chrysene-d12	21.442	240	13181	0.400	ng/ul	0.00
23) Perylene-d12	23.833	264	13320	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.314	96	2262	0.732	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.255	152	7034	0.459	ng/ul	0.00
18) Fluoranthene-d10	19.280	212	16865	0.477	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.348	88	5885	1.743	ng/ul#	87
5) Naphthalene	10.716	128	11993	0.414	ng/ul	100
7) 2-Methylnaphthalene	12.327	142	8259	0.437	ng/ul	99
8) 1-Methylnaphthalene	12.547	142	7651	0.406	ng/ul	99
10) Acenaphthylene	14.224	152	11402	0.486	ng/ul	99
11) Acenaphthene	14.566	153	8200	0.418	ng/ul	99
12) Fluorene	15.551	166	9318	0.415	ng/ul	99
14) Pentachlorophenol	16.904	266	2257	1.128	ng/ul	99
15) Phenanthrene	17.292	178	15011	0.422	ng/ul	99
16) Anthracene	17.381	178	14323	0.466	ng/ul	98
19) Fluoranthene	19.313	202	18841	0.429	ng/ul	98
20) Pyrene	19.675	202	19736	0.444	ng/ul	95
21) Benzo(a)anthracene	21.424	228	21052	0.522	ng/ul	100
22) Chrysene	21.477	228	21539	0.473	ng/ul	99
24) Benzo(b)fluoranthene	23.105	252	22252	0.444	ng/ul	99
25) Benzo(k)fluoranthene	23.151	252	22390	0.444	ng/ul	99
26) Benzo(a)pyrene	23.727	252	19890	0.468	ng/ul	99
27) Indeno(1,2,3-cd)pyrene	26.313	276	25475	0.464	ng/ul#	93
28) Dibenzo(a,h)anthracene	26.330	278	21154	0.467	ng/ul	99
29) Benzo(g,h,i)perylene	27.071	276	21544	0.452	ng/ul	98

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

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