

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN080224\
 Data File : BN033197.D
 Acq On : 02 Aug 2024 09:59
 Operator : MA/JU
 Sample : PB162329BS
 Misc :
 ALS Vial : 31 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS329

Quant Time: Aug 02 10:38:32 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN072524.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Jul 30 14:49:49 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.512	152	2819	0.400	ng/ul	-0.04
4) Naphthalene-d8	10.273	136	9969	0.400	ng/ul	-0.04
9) Acenaphthene-d10	14.138	164	5804	0.400	ng/ul	-0.02
13) Phenanthrene-d10	16.920	188	11676	0.400	ng/ul	-0.02
17) Chrysene-d12	21.134	240	7921	0.400	ng/ul	-0.02
23) Perylene-d12	23.311	264	9656	0.400	ng/ul	-0.03
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.052	96	2730	0.766	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.884	152	5146	0.316	ng/ul	-0.04
18) Fluoranthene-d10	18.958	212	10434	0.465	ng/ul	-0.01
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.081	88	8038	1.977	ng/ul#	66
5) Naphthalene	10.328	128	10255	0.401	ng/ul	98
7) 2-Methylnaphthalene	11.961	142	5813	0.335	ng/ul	98
8) 1-Methylnaphthalene	12.159	142	6042	0.322	ng/ul	98
10) Acenaphthylene	13.866	152	9350	0.394	ng/ul	99
11) Acenaphthene	14.199	153	7437	0.416	ng/ul	99
12) Fluorene	15.221	166	7717	0.371	ng/ul	100
14) Pentachlorophenol	16.616	266	1119	0.471	ng/ul	99
15) Phenanthrene	16.962	178	11185	0.357	ng/ul	99
16) Anthracene	17.063	178	9832	0.349	ng/ul	98
19) Fluoranthene	18.986	202	12901	0.495	ng/ul	100
20) Pyrene	19.358	202	13305	0.476	ng/ul	97
21) Benzo(a)anthracene	21.122	228	9433	0.349	ng/ul	99
22) Chrysene	21.172	228	12334	0.391	ng/ul	99
24) Benzo(b)fluoranthene	22.662	252	10720	0.380	ng/ul	96
25) Benzo(k)fluoranthene	22.703	252	12677	0.375	ng/ul	95
26) Benzo(a)pyrene	23.226	252	11059	0.417	ng/ul#	92
27) Indeno(1,2,3-cd)pyrene	25.500	276	15020	0.372	ng/ul#	100
28) Dibenzo(a,h)anthracene	25.497	278	11648	0.364	ng/ul	99
29) Benzo(g,h,i)perylene	26.181	276	11976	0.373	ng/ul	97

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

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