

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN090623\
 Data File : BN027411.D
 Acq On : 06 Sep 2023 17:37
 Operator : MA/JU
 Sample : PB155237BS
 Misc :
 ALS Vial : 10 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS237

Quant Time: Sep 06 18:15:09 2023
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN090123.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Sat Sep 02 06:16:48 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	8.051	152	5033	0.400	ng/ul	0.00
4) Naphthalene-d8	10.872	136	15121	0.400	ng/ul	-0.01
9) Acenaphthene-d10	14.680	164	9275	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.430	188	22471	0.400	ng/ul	0.00
17) Chrysene-d12	21.606	240	20111	0.400	ng/ul	0.00
23) Perylene-d12	24.123	264	19260	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.410	96	3953	0.590	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.451	152	7582	0.330	ng/ul	0.00
18) Fluoranthene-d10	19.450	212	19704	0.286	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.448	88	10286	1.518	ng/ul#	87
5) Naphthalene	10.922	128	13968	0.318	ng/ul	99
7) 2-Methylnaphthalene	12.522	142	9160	0.334	ng/ul	99
8) 1-Methylnaphthalene	12.742	142	9609	0.329	ng/ul	99
10) Acenaphthylene	14.412	152	12767	0.301	ng/ul	100
11) Acenaphthene	14.745	153	11068	0.303	ng/ul	99
12) Fluorene	15.725	166	12774	0.312	ng/ul	99
14) Pentachlorophenol	17.059	266	3331	0.760	ng/ul	99
15) Phenanthrene	17.473	178	22220	0.304	ng/ul	99
16) Anthracene	17.561	178	18062	0.310	ng/ul	100
19) Fluoranthene	19.478	202	26155	0.283	ng/ul	100
20) Pyrene	19.845	202	26921	0.286	ng/ul	98
21) Benzo(a)anthracene	21.592	228	24228	0.334	ng/ul	99
22) Chrysene	21.647	228	26529	0.314	ng/ul	99
24) Benzo(b)fluoranthene	23.340	252	26837	0.295	ng/ul	99
25) Benzo(k)fluoranthene	23.392	252	26130	0.294	ng/ul	98
26) Benzo(a)pyrene	24.009	252	21713	0.306	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	26.774	276	27249	0.345	ng/ul#	99
28) Dibenzo(a,h)anthracene	26.794	278	21744	0.360	ng/ul	97
29) Benzo(g,h,i)perylene	27.599	276	23288	0.327	ng/ul	99

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

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