

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN092623\
 Data File : BN027874.D
 Acq On : 26 Sep 2023 22:45
 Operator : MA/JU
 Sample : PB155620BS
 Misc :
 ALS Vial : 23 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS620

Quant Time: Sep 27 03:02:16 2023
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN091523.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Sep 22 03:58:00 2023
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.966	152	2650	0.400	ng/ul	0.00
4) Naphthalene-d8	10.784	136	8996	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.606	164	6174	0.400	ng/ul	0.01
13) Phenanthrene-d10	17.358	188	14880	0.400	ng/ul	0.02
17) Chrysene-d12	21.539	240	15224	0.400	ng/ul	0.02
23) Perylene-d12	24.012	264	16710	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.334	96	2606	0.781	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.363	152	6196	0.455	ng/ul	0.00
18) Fluoranthene-d10	19.380	212	18023	0.388	ng/ul	0.02
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.368	88	7714	2.220	ng/ul#	85
5) Naphthalene	10.834	128	9849	0.383	ng/ul	99
7) 2-Methylnaphthalene	12.440	142	6672	0.409	ng/ul	99
8) 1-Methylnaphthalene	12.654	142	6577	0.391	ng/ul	98
10) Acenaphthylene	14.333	152	10342	0.383	ng/ul	99
11) Acenaphthene	14.666	153	8288	0.355	ng/ul	98
12) Fluorene	15.651	166	9835	0.359	ng/ul	99
14) Pentachlorophenol	16.987	266	2942	0.845	ng/ul	98
15) Phenanthrene	17.401	178	16866	0.361	ng/ul	99
16) Anthracene	17.489	178	14711	0.376	ng/ul	99
19) Fluoranthene	19.413	202	20536	0.334	ng/ul	99
20) Pyrene	19.780	202	21390	0.337	ng/ul	99
21) Benzo(a)anthracene	21.521	228	21931	0.372	ng/ul	100
22) Chrysene	21.577	228	22334	0.367	ng/ul	99
24) Benzo(b)fluoranthene	23.243	252	24513	0.329	ng/ul	97
25) Benzo(k)fluoranthene	23.293	252	24763	0.343	ng/ul	98
26) Benzo(a)pyrene	23.898	252	23325	0.389	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	26.609	276	29293	0.382	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.626	278	23914	0.387	ng/ul	98
29) Benzo(g,h,i)perylene	27.417	276	24584	0.377	ng/ul	97

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

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