

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN082824\
 Data File : BN033664.D
 Acq On : 29 Aug 2024 02:49
 Operator : MA/JU
 Sample : PB162960BS
 Misc :
 ALS Vial : 29 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS960

Quant Time: Aug 29 03:21:39 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN082724.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Tue Aug 27 23:48:09 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.536	152	5511	0.400	ng/ul	0.00
4) Naphthalene-d8	10.294	136	14923	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.174	164	7455	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.923	188	15752	0.400	ng/ul	0.00
17) Chrysene-d12	21.130	240	12772	0.400	ng/ul	0.00
23) Perylene-d12	23.292	264	13553	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.143	96	4460	0.787	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.894	152	8301	0.388	ng/ul	0.00
18) Fluoranthene-d10	18.962	212	16045	0.421	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.177	88	11734	1.867	ng/ul#	79
5) Naphthalene	10.343	128	14261	0.357	ng/ul	99
7) 2-Methylnaphthalene	11.966	142	8967	0.355	ng/ul	100
8) 1-Methylnaphthalene	12.191	142	10536	0.405	ng/ul	99
10) Acenaphthylene	13.888	152	11508	0.340	ng/ul	100
11) Acenaphthene	14.235	153	9322	0.381	ng/ul	99
12) Fluorene	15.229	166	10149	0.372	ng/ul	100
14) Pentachlorophenol	16.581	266	2948	0.595	ng/ul	99
15) Phenanthrene	16.965	178	16296	0.351	ng/ul	100
16) Anthracene	17.054	178	14532	0.348	ng/ul	100
19) Fluoranthene	18.990	202	18455	0.366	ng/ul	98
20) Pyrene	19.352	202	19193	0.359	ng/ul	99
21) Benzo(a)anthracene	21.115	228	17655	0.351	ng/ul	99
22) Chrysene	21.168	228	18258	0.359	ng/ul	99
24) Benzo(b)fluoranthene	22.649	252	18429	0.350	ng/ul	99
25) Benzo(k)fluoranthene	22.690	252	19242	0.366	ng/ul	98
26) Benzo(a)pyrene	23.196	252	16763	0.397	ng/ul	98
27) Indeno(1,2,3-cd)pyrene	25.438	276	22613	0.356	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.452	278	17575	0.357	ng/ul	99
29) Benzo(g,h,i)perylene	26.089	276	17805	0.359	ng/ul	100

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

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