

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033957.D
 Acq On : 15 Sep 2024 18:15
 Operator : JU/RC
 Sample : PB163312BS
 Misc :
 ALS Vial : 48 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS312

Quant Time: Sep 16 00:14:22 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 : Sun Sep 15 01:52:43 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.481	152	3985	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	11071	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.128	164	5661	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.881	188	11856	0.400	ng/ul	0.00
17) Chrysene-d12	21.095	240	9975	0.400	ng/ul	0.00
23) Perylene-d12	23.237	264	10923	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	3323	0.811	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	6154	0.387	ng/ul	0.00
18) Fluoranthene-d10	18.920	212	13161	0.443	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	8755	1.926	ng/ul#	80
5) Naphthalene	10.288	128	10352	0.350	ng/ul	99
7) 2-Methylnaphthalene	11.916	142	6465	0.345	ng/ul	100
8) 1-Methylnaphthalene	12.136	142	7786	0.404	ng/ul	99
10) Acenaphthylene	13.841	152	8430	0.328	ng/ul	99
11) Acenaphthene	14.188	153	6949	0.374	ng/ul	98
12) Fluorene	15.183	166	7522	0.363	ng/ul	98
14) Pentachlorophenol	16.539	266	657	0.176	ng/ul	98
15) Phenanthrene	16.923	178	12168	0.348	ng/ul	99
16) Anthracene	17.016	178	10484	0.334	ng/ul	100
19) Fluoranthene	18.952	202	14595	0.371	ng/ul	99
20) Pyrene	19.315	202	15359	0.368	ng/ul	99
21) Benzo(a)anthracene	21.080	228	13269	0.338	ng/ul	99
22) Chrysene	21.130	228	15587	0.393	ng/ul	100
24) Benzo(b)fluoranthene	22.602	252	14738	0.348	ng/ul	98
25) Benzo(k)fluoranthene	22.643	252	16675	0.393	ng/ul	99
26) Benzo(a)pyrene	23.143	252	13789	0.405	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.348	276	18700	0.366	ng/ul	99
28) Dibenzo(a,h)anthracene	25.368	278	14460	0.364	ng/ul	97
29) Benzo(g,h,i)perylene	25.992	276	15398	0.385	ng/ul	98

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

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