

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033958.D
 Acq On : 15 Sep 2024 18:51
 Operator : JU/RC
 Sample : PB163357BS
 Misc :
 ALS Vial : 49 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS357

Quant Time: Sep 16 00:14:34 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	3555	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	9821	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.123	164	4963	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.881	188	10299	0.400	ng/ul	0.00
17) Chrysene-d12	21.095	240	8615	0.400	ng/ul	0.00
23) Perylene-d12	23.234	264	9551	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	3779	1.034	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	6784	0.481	ng/ul	0.00
18) Fluoranthene-d10	18.920	212	13391	0.521	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.131	88	9249	2.281	ng/ul#	79
5) Naphthalene	10.288	128	10972	0.418	ng/ul	99
7) 2-Methylnaphthalene	11.916	142	6753	0.406	ng/ul	98
8) 1-Methylnaphthalene	12.136	142	8094	0.473	ng/ul	99
10) Acenaphthylene	13.841	152	8743	0.388	ng/ul	99
11) Acenaphthene	14.188	153	7167	0.441	ng/ul	98
12) Fluorene	15.183	166	7709	0.424	ng/ul	100
14) Pentachlorophenol	16.543	266	719	0.222	ng/ul	99
15) Phenanthrene	16.923	178	12065	0.397	ng/ul	100
16) Anthracene	17.016	178	10388	0.380	ng/ul	100
19) Fluoranthene	18.952	202	13911	0.409	ng/ul	100
20) Pyrene	19.315	202	14583	0.404	ng/ul	99
21) Benzo(a)anthracene	21.080	228	12671	0.374	ng/ul	99
22) Chrysene	21.130	228	15163	0.443	ng/ul	99
24) Benzo(b)fluoranthene	22.602	252	14510	0.391	ng/ul	98
25) Benzo(k)fluoranthene	22.646	252	16224	0.437	ng/ul	98
26) Benzo(a)pyrene	23.146	252	13243	0.445	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.351	276	18467	0.413	ng/ul#	98
28) Dibenzo(a,h)anthracene	25.368	278	14226	0.410	ng/ul	97
29) Benzo(g,h,i)perylene	25.992	276	15192	0.435	ng/ul	99

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

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