

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN091424\
 Data File : BN033942.D
 Acq On : 15 Sep 2024 08:02
 Operator : JU/RC
 Sample : PB163350BS
 Misc :
 ALS Vial : 32 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS350

Quant Time: Sep 16 00:12:25 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	7.485	152	3664	0.400	ng/ul	0.00
4) Naphthalene-d8	10.239	136	10131	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.128	164	5180	0.400	ng/ul	0.00
13) Phenanthrene-d10	16.881	188	10838	0.400	ng/ul	0.00
17) Chrysene-d12	21.098	240	9449	0.400	ng/ul	0.00
23) Perylene-d12	23.240	264	10404	0.400	ng/ul	0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.097	96	3586	0.952	ng/ul	0.00
6) 2-Methylnaphthalene-d10	11.839	152	6460	0.444	ng/ul	0.00
18) Fluoranthene-d10	18.925	212	13224	0.469	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.131	88	8726	2.088	ng/ul#	81
5) Naphthalene	10.288	128	10233	0.378	ng/ul	99
7) 2-Methylnaphthalene	11.916	142	6449	0.376	ng/ul	100
8) 1-Methylnaphthalene	12.136	142	7707	0.437	ng/ul	99
10) Acenaphthylene	13.841	152	8322	0.354	ng/ul	99
11) Acenaphthene	14.188	153	6836	0.403	ng/ul	99
12) Fluorene	15.188	166	7362	0.388	ng/ul	99
14) Pentachlorophenol	16.543	266	705	0.207	ng/ul	98
15) Phenanthrene	16.923	178	11665	0.365	ng/ul	99
16) Anthracene	17.016	178	10246	0.357	ng/ul	99
19) Fluoranthene	18.952	202	13730	0.368	ng/ul	99
20) Pyrene	19.320	202	14409	0.364	ng/ul	98
21) Benzo(a)anthracene	21.083	228	12963	0.348	ng/ul	99
22) Chrysene	21.133	228	14630	0.389	ng/ul	99
24) Benzo(b)fluoranthene	22.605	252	14726	0.365	ng/ul	98
25) Benzo(k)fluoranthene	22.646	252	15595	0.386	ng/ul	99
26) Benzo(a)pyrene	23.146	252	13283	0.410	ng/ul	97
27) Indeno(1,2,3-cd)pyrene	25.354	276	18211	0.374	ng/ul#	99
28) Dibenzo(a,h)anthracene	25.371	278	14113	0.373	ng/ul	97
29) Benzo(g,h,i)perylene	25.998	276	14791	0.389	ng/ul	99

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

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