

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN030224\
 Data File : BN030243.D
 Acq On : 04 Mar 2024 06:48
 Operator : MA/JU
 Sample : PB159151BS
 Misc :
 ALS Vial : 66 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS151

Quant Time: Mar 04 07:16:51 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN030224.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Mon Mar 04 05:54:20 2024
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.834	152	1231	0.400	ng/ul	0.00
4) Naphthalene-d8	10.632	136	3345	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.482	164	1563	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.242	188	2920	0.400	ng/ul	0.00
17) Chrysene-d12	21.445	240	2700	0.400	ng/ul	# 0.00
23) Perylene-d12	23.798	264	3131	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.307	96	1398	0.820	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.233	152	1645	0.380	ng/ul	0.00
18) Fluoranthene-d10	19.276	212	3317	0.377	ng/ul	0.00
Target Compounds						
						Qvalue
2) 1,4-Dioxane	3.336	88	4334	2.344	ng/ul#	86
5) Naphthalene	10.682	128	3794	0.384	ng/ul	98
7) 2-Methylnaphthalene	12.310	142	2012	0.376	ng/ul	98
8) 1-Methylnaphthalene	12.519	142	2099	0.370	ng/ul	99
10) Acenaphthylene	14.200	152	3387	0.388	ng/ul	97
11) Acenaphthene	14.542	153	2432	0.379	ng/ul	99
12) Fluorene	15.541	166	2513	0.385	ng/ul	98
14) Pentachlorophenol	16.887	266	1013	1.216	ng/ul	97
15) Phenanthrene	17.280	178	3753	0.383	ng/ul	98
16) Anthracene	17.377	178	3334	0.387	ng/ul	100
19) Fluoranthene	19.304	202	4464	0.372	ng/ul	99
20) Pyrene	19.666	202	4865	0.373	ng/ul	98
21) Benzo(a)anthracene	21.430	228	4052	0.425	ng/ul	99
22) Chrysene	21.483	228	4596	0.401	ng/ul	99
24) Benzo(b)fluoranthene	23.084	252	4923	0.408	ng/ul	97
25) Benzo(k)fluoranthene	23.134	252	5586	0.392	ng/ul	96
26) Benzo(a)pyrene	23.696	252	4875	0.397	ng/ul#	93
27) Indeno(1,2,3-cd)pyrene	26.226	276	6405	0.408	ng/ul#	97
28) Dibenzo(a,h)anthracene	26.260	278	5058	0.416	ng/ul	94
29) Benzo(g,h,i)perylene	26.967	276	5696	0.391	ng/ul	95

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

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