

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN111222\
 Data File : BN022622.D
 Acq On : 11 Nov 2022 21:27
 Operator : CG/JU
 Sample : PB148943BS
 Misc :
 ALS Vial : 14 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS943

Quant Time: Nov 11 22:41:16 2022
 Quant Method : Z:\SVOASRV\HPCHEM1\BNA_N\METHODS\SFAM-EPA-SIM-BN102022.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Fri Nov 11 22:23:04 2022
 Response via : Initial Calibration

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

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.902	152	1759	0.400	ng/ul	0.00
4) Naphthalene-d8	10.723	136	5462	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.577	164	3035	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.336	188	5851	0.400	ng/ul	0.00
17) Chrysene-d12	21.544	240	4962	0.400	ng/ul	0.00
23) Perylene-d12	23.985	264	4003	0.400	ng/ul	# 0.00
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.211	96	1867	0.817	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.323	152	3577	0.433	ng/ul	0.00
18) Fluoranthene-d10	19.375	212	6852	0.411	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.245	88	5434	2.315	ng/ul#	88
5) Naphthalene	10.772	128	6555	0.406	ng/ul	98
7) 2-Methylnaphthalene	12.400	142	4139	0.407	ng/ul	100
8) 1-Methylnaphthalene	12.615	142	4327	0.414	ng/ul	99
10) Acenaphthylene	14.299	152	5532	0.406	ng/ul	97
11) Acenaphthene	14.642	153	4496	0.391	ng/ul	99
12) Fluorene	15.632	166	4999	0.375	ng/ul	98
14) Pentachlorophenol	16.994	266	1058	0.761	ng/ul	99
15) Phenanthrene	17.379	178	7776	0.396	ng/ul	98
16) Anthracene	17.471	178	6665	0.403	ng/ul	97
19) Fluoranthene	19.407	202	8627	0.380	ng/ul	96
20) Pyrene	19.770	202	8711	0.382	ng/ul	99
21) Benzo(a)anthracene	21.529	228	7665	0.412	ng/ul	98
22) Chrysene	21.582	228	8306	0.421	ng/ul	99
24) Benzo(b)fluoranthene	23.233	252	7955	0.396	ng/ul	96
25) Benzo(k)fluoranthene	23.283	252	7529	0.385	ng/ul	94
26) Benzo(a)pyrene	23.877	252	7255	0.453	ng/ul	95
27) Indeno(1,2,3-cd)pyrene	26.541	276	7832	0.393	ng/ul#	94
28) Dibenzo(a,h)anthracene	26.565	278	6463	0.406	ng/ul#	93
29) Benzo(g,h,i)perylene	27.336	276	6760	0.417	ng/ul	97

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

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