

Data Path : Z:\svoasrv\HPCHEM1\BNA_N\Data\BN020824\
 Data File : BN029598.D
 Acq On : 09 Feb 2024 03:59
 Operator : MA/JU
 Sample : PB158770BS
 Misc :
 ALS Vial : 26 Sample Multiplier: 1

Instrument :
 BNA_N
 ClientSampleId :
 SLCS770

Quant Time: Feb 09 04:30:10 2024
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_N\Methods\SFAM-EPA-SIM-BN020624.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Wed Feb 07 02:17:13 2024
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)
Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.888	152	5242	0.400	ng/ul	0.00
4) Naphthalene-d8	10.682	136	14548	0.400	ng/ul	0.00
9) Acenaphthene-d10	14.528	164	6433	0.400	ng/ul	0.00
13) Phenanthrene-d10	17.279	188	11910	0.400	ng/ul	0.00
17) Chrysene-d12	21.473	240	8863	0.400	ng/ul	0.00
23) Perylene-d12	23.847	264	9090	0.400	ng/ul	-0.01
System Monitoring Compounds						
3) 1,4-Dioxane-d8	3.336	96	4850	0.742	ng/ul	0.00
6) 2-Methylnaphthalene-d10	12.277	152	7209	0.389	ng/ul	0.00
18) Fluoranthene-d10	19.307	212	11273	0.387	ng/ul	0.00
Target Compounds						
					Qvalue	
2) 1,4-Dioxane	3.366	88	15872	2.231	ng/ul#	87
5) Naphthalene	10.731	128	16358	0.400	ng/ul	99
7) 2-Methylnaphthalene	12.348	142	9325	0.397	ng/ul	100
8) 1-Methylnaphthalene	12.568	142	9269	0.388	ng/ul	100
10) Acenaphthylene	14.246	152	12542	0.400	ng/ul	99
11) Acenaphthene	14.588	153	9510	0.399	ng/ul	99
12) Fluorene	15.578	166	9837	0.394	ng/ul	100
14) Pentachlorophenol	16.920	266	2434	0.916	ng/ul#	51
15) Phenanthrene	17.321	178	14953	0.402	ng/ul	100
16) Anthracene	17.410	178	12811	0.409	ng/ul	99
19) Fluoranthene	19.340	202	15867	0.390	ng/ul	99
20) Pyrene	19.702	202	16535	0.396	ng/ul	98
21) Benzo(a)anthracene	21.459	228	13815	0.432	ng/ul	99
22) Chrysene	21.511	228	14987	0.429	ng/ul	100
24) Benzo(b)fluoranthene	23.124	252	14595	0.427	ng/ul	99
25) Benzo(k)fluoranthene	23.174	252	16179	0.427	ng/ul	98
26) Benzo(a)pyrene	23.741	252	13848	0.431	ng/ul	100
27) Indeno(1,2,3-cd)pyrene	26.302	276	17133	0.439	ng/ul#	100
28) Dibenzo(a,h)anthracene	26.333	278	13431	0.439	ng/ul	99
29) Benzo(g,h,i)perylene	27.050	276	15182	0.428	ng/ul	98

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

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